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# DP-GEN

## DeepModeling

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# OVERVIEW

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## OVERVIEW

### 1.1 About DP-GEN

DP-GEN (Deep Generator) is a software written in Python, delicately designed to generate a deep learning based model of interatomic potential energy and force field. DP-GEN is dependent on [DeepMD-kit](#). With highly scalable interface with common softwares for molecular simulation, DP-GEN is capable to automatically prepare scripts and maintain job queues on HPC machines (High Performance Cluster) and analyze results.

If you use this software in any publication, please cite:

Yuzhi Zhang, Haidi Wang, Weijie Chen, Jinzhe Zeng, Linfeng Zhang, Han Wang, and Weinan E, DP-GEN: A concurrent learning platform for the generation of reliable deep learning based potential energy models, *Computer Physics Communications*, 2020, 107206.

#### 1.1.1 Highlighted features

- **Accurate and efficient:** DP-GEN is capable to sample more than tens of million structures and select only a few for first principles calculation. DP-GEN will finally obtain a uniformly accurate model.
- **User-friendly and automatic:** Users may install and run DP-GEN easily. Once succusefully running, DP-GEN can dispatch and handle all jobs on HPCs, and thus there's no need for any personal effort.
- **Highly scalable:** With modularized code structures, users and developers can easily extend DP-GEN for their most relevant needs. DP-GEN currently supports for HPC systems (Slurm, PBS, LSF and cloud machines ), Deep Potential interface with DeePMD-kit, MD interface with [LAMMPS](#), [Gromacs](#) and *ab-initio* calculation interface with VASP, PWSCF, CP2K, SIESTA and Gaussian, Abacus, PWMAT, etc . We're sincerely welcome and embraced to users' contributions, with more possibilities and cases to use DP-GEN.

### 1.2 Download and install

DP-GEN only supports Python 3.9 and above. You can use one of the following methods to install DP-GEN:

- Install via pip: `pip install dpgen`
- Install via conda: ``conda install -c conda-forge dpgen``
- Install from source code: `git clone https://github.com/deepmodeling/dpgen && pip install ./dpgen`

To test if the installation is successful, you may execute

```
dpgen -h
```

## 1.3 Use DP-GEN

A quick-start on using DPGEN can be found [here](#). You can follow the [Handson tutorial](#), it is friendly to new users.

## 1.4 Case Studies

- [Practical-Guidelines-for-DP](#)

Before starting a new Deep Potential (DP) project, we suggest people (especially those who are newbies) read the following context first to get some insights into what tools we can use, what kinds of risks and difficulties we may meet, and how we can advance a new DP project smoothly.

- [Convergence-Test](#)

to ensure the data quality, the reliability of the final model, as well as the feasibility of the project, a convergence test should be done first.

- [Gas-phase](#)

In this tutorial, we will take the simulation of methane combustion as an example and introduce the procedure of DP-based MD simulation.

- [Mg-Y\\_alloy](#)

We will briefly analyze the candidate configurational space of a metallic system by taking Mg-based Mg-Y binary alloy as an example. The task is divided into steps during the DP-GEN process.

- [Transfer-learning](#)

This tutorial will introduce how to implement potential energy surface (PES) transfer-learning by using the DP-GEN software. In DP-GEN (version > 0.8.0), the “simplify” module is designed for this purpose.

## 1.5 License

The project dpgen is licensed under [GNU LGPLv3.0](#)

## COMMAND LINE INTERFACE

dpngen is a convenient script that uses DeepGenerator to prepare initial data, drive DeepMDkit and analyze results. This script works based on several sub-commands with their own options. To see the options for the sub-commands, type “dpngen sub-command -h”.

```
usage: dpngen [-h]
              {init_surf,init_bulk,auto_gen_param,init_reaction,run,run/report,collect,
              ↪simplify,autotest,db,gui}
              ...
```

### 2.1 Sub-commands

#### 2.1.1 init\_surf

Generating initial data for surface systems.

```
dpngen init_surf [-h] PARAM [MACHINE]
```

##### Positional Arguments

<b>PARAM</b>	parameter file, json/yaml format
<b>MACHINE</b>	machine file, json/yaml format

#### 2.1.2 init\_bulk

Generating initial data for bulk systems.

```
dpngen init_bulk [-h] PARAM [MACHINE]
```

### Positional Arguments

<b>PARAM</b>	parameter file, json/yaml format
<b>MACHINE</b>	machine file, json/yaml format

#### 2.1.3 auto\_gen\_param

auto gen param.json

```
dpngen auto_gen_param [-h] PARAM
```

### Positional Arguments

<b>PARAM</b>	parameter file, json/yaml format
--------------	----------------------------------

#### 2.1.4 init\_reaction

Generating initial data for reactive systems.

```
dpngen init_reaction [-h] PARAM [MACHINE]
```

### Positional Arguments

<b>PARAM</b>	parameter file, json/yaml format
<b>MACHINE</b>	machine file, json/yaml format

#### 2.1.5 run

Main process of Deep Potential Generator.

```
dpngen run [-h] [-d] PARAM MACHINE
```

### Positional Arguments

<b>PARAM</b>	parameter file, json/yaml format
<b>MACHINE</b>	machine file, json/yaml format



## Named Arguments

**-d, --debug**      log debug info  
Default: False

### 2.1.6 run/report

Report the systems and the thermodynamic conditions of the labeled frames.

```
dpgen run/report [-h] [-s] [-i] [-t] [-p PARAM] [-v] JOB_DIR
```

## Positional Arguments

**JOB\_DIR**      the directory of the DP-GEN job,

## Named Arguments

**-s, --stat-sys**      count the labeled frames for each system  
Default: False

**-i, --stat-iter**      print the iteration candidate,failed,accurate count and fp calculation,success and fail count  
Default: False

**-t, --stat-time**      print the iteration time, warning!! assume model\_devi parallel cores == 1  
Default: False

**-p, --param**      the json file provides DP-GEN paramters, should be located in JOB\_DIR  
Default: "param.json"

**-v, --verbose**      being loud  
Default: False

### 2.1.7 collect

Collect data.

```
dpgen collect [-h] [-p PARAMETER] [-v] [-m] [-s] JOB_DIR OUTPUT
```

## Positional Arguments

**JOB\_DIR**      the directory of the DP-GEN job  
**OUTPUT**      the output directory of data

### Named Arguments

<b>-p, --parameter</b>	the json file provides DP-GEN paramters, should be located in JOB_DIR Default: "param.json"
<b>-v, --verbose</b>	print number of data in each system Default: False
<b>-m, --merge</b>	merge the systems with the same chemical formula Default: False
<b>-s, --shuffle</b>	shuffle the data systems Default: False

### 2.1.8 simplify

Simplify data.

```
dpngen simplify [-h] [-d] PARAM MACHINE
```

### Positional Arguments

<b>PARAM</b>	parameter file, json/yaml format
<b>MACHINE</b>	machine file, json/yaml format

### Named Arguments

<b>-d, --debug</b>	log debug info Default: False
--------------------	----------------------------------

### 2.1.9 autotest

Auto-test for Deep Potential.

```
dpngen autotest [-h] [-d] TASK PARAM [MACHINE]
```

### Positional Arguments

<b>TASK</b>	task can be make, run or post
<b>PARAM</b>	parameter file, json/yaml format
<b>MACHINE</b>	machine file, json/yaml format

### Named Arguments

<b>-d, --debug</b>	log debug info
	Default: False

#### 2.1.10 db

Collecting data from DP-GEN.

```
dpgen db [-h] PARAM
```

### Positional Arguments

<b>PARAM</b>	parameter file, json format
--------------	-----------------------------

#### 2.1.11 gui

Serve DP-GUI.

```
dpgen gui [-h] [-p PORT] [--bind_all]
```

### Named Arguments

<b>-p, --port</b>	The port to serve DP-GUI on. Default: 6042
<b>--bind_all</b>	Serve on all public interfaces. This will expose your DP-GUI instance to the network on both IPv4 and IPv6 (where available). Default: False



## CODE STRUCTURE

Let's look at the home page of DP-GEN. <https://github.com/deepmodeling/dpgen>

```
build
CITATION.cff
conda
dist
doc
dpgen
dpgen.egg-info
examples
LICENSE
README.md
requirements.txt
setup.py
tests
```


- `tests` : unittest tools for developers.
- `examples`: templates for PARAM and MACHINE files for different software, versions and tasks. For details of the parameters in PARAM, you can refer to TASK parameters chapters in this document. If you are confused about how to set up a JSON file, you can also use `dpgui`

Most of the code related to DP-GEN functions is in the `dpgen` directory. Open the `dpgen` directory, and we can see

```
arginfo.py
auto_test
collect
data
database
_date.py
dispatcher
generator
__init__.py
main.py
__pycache__
remote
simplify
tools
util.py
_version.py
```

- `auto_test` corresponds to `dpgen autotest`, for undertaking materials property analysis.

- `collect` corresponds to `dpgen collect`.
- `data` corresponds to `dpgen init_bulk`, `dpgen init_surf` and `dpgen init_reaction`, for preparing initial data of bulk and surf systems.
- `database` is the source code for collecting data generated by DP-GEN and interface with database.
- `simplify` corresponds to `dpgen simplify`.
- `remote` and `dispatcher` : source code for automatically submitting scripts, maintaining job queues and collecting results. **Notice this part has been integrated into `dpdispatcher`** generator is the core part of DP-GEN. It's for main process of deep generator. Let's open this folder.



```
arginfo.py
ch4
__init__.py
lib
run.py
```

`run.py` is the core of DP-GEN, corresponding to `dpgen run`. We can find `make_train`, `run_train`, ... `post_fp`, and other steps related functions here.

## 4.1 Overview of the Run process

The run process contains a series of successive iterations, undertaken in order such as heating the system to certain temperatures. Each iteration is composed of three steps: exploration, labeling, and training. Accordingly, there are three sub-folders: 00.train, 01.model\_devi, and 02.fp in each iteration.

00.train: DP-GEN will train several (default 4) models based on initial and generated data. The only difference between these models is the random seed for neural network initialization.

01.model\_devi : represent for model-deviation. DP-GEN will use models obtained from 00.train to run Molecular Dynamics(default LAMMPS). Larger deviation for structure properties (default is the force of atoms) means less accuracy of the models. Using this criterion, a few structures will be selected and put into the next stage 02.fp for more accurate calculation based on First Principles.

02.fp : Selected structures will be calculated by first-principles methods(default VASP). DP-GEN will obtain some new data and put them together with initial data and data generated in previous iterations. After that, new training will be set up and DP-GEN will enter the next iteration!

In the run process of the DP-GEN, we need to specify the basic information about the system, the initial data, and details of the training, exploration, and labeling tasks. In addition, we need to specify the software, machine environment, and computing resource and enable the process of job generation, submission, query, and collection automatically. We can perform the run process as we expect by specifying the keywords in param.json and machine.json, and they will be introduced in detail in the following sections.

Here, we give a general description of the run process. We can execute the run process of DP-GEN easily by:

```
dpgen run param.json machine.json
```

The following files or folders will be created and upgraded by codes

- iter.00000x contains the main results that DP-GEN generates in the first iteration.
- record.dpgen records the current stage of the run process.
- dpgen.log includes time and iteration information.

When the first iteration is completed, the folder structure of iter.000000 is like this:

```
$ ls iter.000000
00.train 01.model_devi 02.fp
```

In folder iter.000000/ 00.train:

- Folder 00x contains the input and output files of the DeePMD-kit, in which a model is trained.

- graph.00x.pb is the model DeePMD-kit generates. The only difference between these models is the random seed for neural network initialization.

In folder iter.000000/ 01.model\_devi

- Folder confs contains the initial configurations for LAMMPS MD converted from POSCAR you set in *sys\_confs* of param.json.
- Folder task.000.00000x contains the input and output files of the LAMMPS. In folder task.000.00000x, file model\_devi.out records the model deviation of concerned labels, energy and force in MD. It serves as the criterion for selecting which structures and doing first-principle calculations.

In folder iter.000000/ 02.fp

- candidate.shuffle.000.out records which structures will be selected from last step 01.model\_devi. There are always far more candidates than the maximum you expect to calculate at one time. In this condition, DP-GEN will randomly choose up to *fp\_task\_max* structures and form the folder task.\*.
- rest\_accurate.shuffle.000.out records the other structures where our model is accurate (*max\_devi\_f* is less than *model\_devi\_f\_trust\_lo*, no need to calculate any more),
- rest\_failed.shuffled.000.out records the other structures where our model is too inaccurate (larger than *model\_devi\_f\_trust\_hi*, there may be some error).
- data.000: After first-principle calculations, DP-GEN will collect these data and change them into the format DeePMD-kit needs. In the next iteration's 00.train, these data will be trained together as well as the initial data.

DP-GEN identifies the stage of the run process by a record file, record.dpgen, which will be created and upgraded by codes. Each line contains two numbers: the first is the index of iteration, and the second, ranging from 0 to 9, records which stage in each iteration is currently running.

Index of iterations	Stage in each iteration	Process
0	0	make_train
0	1	run_train
0	2	post_train
0	3	make_model_devi
0	4	run_model_devi
0	5	post_model_devi
0	6	make_fp
0	7	run_fp
0	8	post_fp

0,1,2 correspond to make\_train, run\_train, post\_train. DP-GEN will write scripts in make\_train, run the task by specific machine in run\_train and collect result in post\_train. The records for model\_devi and fp stage follow similar rules.

If the process of DP-GEN stops for some reasons, DP-GEN will automatically recover the main process by record.dpgen. You may also change it manually for your purpose, such as removing the last iterations and recovering from one checkpoint. When re-running dpgen, the process will start from the stage that the last line record.



## 4.2 Example-of-param.json

We have provided different examples of param.json in dpgen/examples/run/. In this section, we give a description of the param.json, taking dpgen/examples/run/dp2.x-lammps-vasp/param\_CH4\_deepmd-kit-2.0.1.json as an example. This is a param.json for a gas-phase methane molecule. Here, DeePMD-kit (v2.x), LAMMPS and VASP codes are used for training, exploration and labeling respectively.

### 4.2.1 basics

The basics related keys in param.json are given as follows

```
"type_map": [
  "H",
  "C"
],
"mass_map": [
  1,
  12
],
```

The basics related keys specify the basic information about the system. *type\_map* gives the atom types, i.e. “H” and “C”. *mass\_map* gives the standard atom weights, i.e. “1” and “12”.

### 4.2.2 data

The data related keys in param.json are given as follows

```
"init_data_prefix": "...../init/",
"init_data_sys": [
  "CH4.POSCAR.01x01x01/02.md/sys-0004-0001/deepmd"
],

"sys_configs_prefix": "...../init/",
"sys_configs": [
  [
    "CH4.POSCAR.01x01x01/01.scale_pert/sys-0004-0001/scale*/00000*/POSCAR"
  ],
  [
    "CH4.POSCAR.01x01x01/01.scale_pert/sys-0004-0001/scale*/00001*/POSCAR"
  ]
],
```

The data related keys specify the init data for training initial DP models and structures used for model\_devi calculations. *init\_data\_prefix* and *init\_data\_sys* specify the location of the init data. *sys\_configs\_prefix* and *sys\_configs* specify the location of the structures.

Here, the init data is provided at “..... /init/CH4.POSCAR.01x01x01/02.md/sys-0004-0001/deepmd”. These structures are divided into two groups and provided at “...../init/CH4.POSCAR.01x01x01/01.scale\_pert/sys-0004-0001/scale\*/00000\*/POSCAR” and “...../init/CH4.POSCAR.01x01x01/01.scale\_pert/sys-0004-0001/scale\*/00001\*/POSCAR”.

### 4.2.3 training

The training related keys in param.json are given as follows

```
"numb_models": 4,
"default_training_param": {
  },
```

The training related keys specify the details of training tasks. `numb_models` specifies the number of models to be trained. “default\_training\_param” specifies the training parameters for deepmd-kit.

Here, 4 DP models will be trained in `00.train`. A detailed explanation of training parameters can be found in DeePMD-kit’s documentation (<https://docs.deepmodeling.com/projects/deepmd/en/master/>).

### 4.2.4 exploration

The exploration related keys in param.json are given as follows

```
"model_devi_dt": 0.002,
"model_devi_skip": 0,
"model_devi_f_trust_lo": 0.05,
"model_devi_f_trust_hi": 0.15,
"model_devi_clean_traj": true,
"model_devi_jobs": [
  {
    "sys_idx": [
      0
    ],
    "temps": [
      100
    ],
    "press": [
      1.0
    ],
    "trj_freq": 10,
    "nsteps": 300,
    "ensemble": "nvt",
    "_idx": "00"
  },
  {
    "sys_idx": [
      1
    ],
    "temps": [
      100
    ],
    "press": [
      1.0
    ],
    "trj_freq": 10,
    "nsteps": 3000,
    "ensemble": "nvt",
    "_idx": "01"
```

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```
}
],
```

The exploration related keys specify the details of exploration tasks. `model_devi_dt` specifies timestep for MD simulation. `model_devi_skip` specifies the number of structures skipped for saving in each MD. `model_devi_f_trust_lo` and `model_devi_f_trust_hi` specify the lower and upper bound of model\_devi of forces for the selection. `model_devi_clean_traj` specifies whether to clean traj folders in MD. If type of model\_devi\_clean\_traj is boolean type then it denote whether to clean traj folders in MD since they are too large. In `model_devi_jobs`, `sys_idx` specifies the group of structures used for model\_devi calculations, `temps` specifies the temperature (K) in MD, `press` specifies the pressure (Bar) in MD, `trj_freq` specifies the frequency of trajectory saved in MD, `nsteps` specifies the running steps of MD, `ensemble` specifies the ensemble used in MD, and “\_idx” specifies the index of iteration.

Here, MD simulations are performed at the temperature of 100 K and the pressure of 1.0 Bar with an integrator time of 2 fs under the nvt ensemble. Two iterations are set in `model_devi_jobs`. MD simulations are run for 300 and 3000 time steps with the first and second groups of structures in `sys_configs` in 00 and 01 iterations. We choose to save all structures generated in MD simulations and have set `trj_freq` as 10, so 30 and 300 structures are saved in 00 and 01 iterations. If the “max\_devi\_f” of saved structure falls between 0.05 and 0.15, DP-GEN will treat the structure as a candidate. We choose to clean traj folders in MD since they are too large. If you want to save the most recent n iterations of traj folders, you can set `model_devi_clean_traj` to be an integer.

## 4.2.5 labeling

The labeling related keys in param.json are given as follows

```
"fp_style": "vasp",
"shuffle_poscar": false,
"fp_task_max": 20,
"fp_task_min": 1,
"fp_pp_path": "...../methane/",
"fp_pp_files": [
  "POTCAR"
],
"fp_incar": "...../INCAR_methane"
```

The labeling related keys specify the details of labeling tasks. `fp_style` specifies software for First Principles. `fp_task_max` and `fp_task_min` specify the minimum and maximum of structures to be calculated in 02.fp of each iteration. `fp_pp_path` and `fp_pp_files` specify the location of the psuedo-potential file to be used for 02.fp. `run_jdata[fp_style=vasp]/fp_incar` specifies input file for VASP. INCAR must specify KSPACING and KGAMMA.

Here, a minimum of 1 and a maximum of 20 structures will be labeled using the VASP code with the INCAR provided at “...../INCAR\_methane” and POTCAR provided at “...../methane/POTCAR” in each iteration. Note that the order of elements in POTCAR should correspond to the order in `type_map`.

All the keys of the DP-GEN are explained in detail in the section Parameters.

## 4.3 Example of machine.json

### 4.3.1 DPDispatcher Update Note

DPDispatcher has updated and the api of machine.json is changed. DP-GEN will use the new DPDispatcher if the value of key `api_version` in machine.json is equal to or large than 1.0. And for now, DPDispatcher is maintained on a separate repo (<https://github.com/deepmodeling/dpdispatcher>). Please check the documents (<https://deepmd.readthedocs.io/projects/dpdispatcher/en/latest/>) for more information about the new DPDispatcher.

DP-GEN will use the old DPDispatcher if the key `api_version` is not specified in machine.json or the `api_version` is smaller than 1.0. This gurantees that the old machine.json still works.

### 4.3.2 New DPDispatcher

Each iteration in the run process of DP-GEN is composed of three steps: exploration, labeling, and training. Accordingly, machine.json is composed of three parts: train, model\_devi, and fp. Each part is a list of dicts. Each dict can be considered as an independent environment for calculation.

In this section, we will show you how to perform train task at a local workstation, model\_devi task at a local Slurm cluster, and fp task at a remote PBS cluster using the new DPDispatcher. For each task, three types of keys are needed:

- Command: provides the command used to execute each step.
- Machine: specifies the machine environment (local workstation, local or remote cluster, or cloud server).
- Resources: specify the number of groups, nodes, CPU, and GPU; enable the virtual environment.

#### Performing train task at a local workstation

In this example, we perform the train task on a local workstation.

```
"train":
{
  "command": "dp",
  "machine": {
    "batch_type": "Shell",
    "context_type": "local",
    "local_root": "./",
    "remote_root": "/home/user1234/work_path"
  },
  "resources": {
    "number_node": 1,
    "cpu_per_node": 4,
    "gpu_per_node": 1,
    "group_size": 1,
    "source_list": ["/home/user1234/deepmd.env"]
  }
},
```

The `command` for the train task in the DeepMD-kit is “dp”.

In machine parameters, `batch_type` specifies the type of job scheduling system. If there is no job scheduling system, we can use the “Shell” to perform the task. `context_type` specifies the method of data transfer, and “local” means copying and moving data via local file storage systems (e.g. cp, mv, etc.). In DP-GEN, the paths of all tasks are

automatically located and set by the software, and therefore `local_root` is always set to `“.”`. The input file for each task will be sent to the `remote_root` and the task will be performed there, so we need to make sure that the path exists.

In the resources parameter, `number_node`, `cpu_per_node`, and `gpu_per_node` specify the number of nodes, the number of CPUs, and the number of GPUs required for a task respectively. `group_size`, which needs to be highlighted, specifies how many tasks will be packed into a group. In the training tasks, we need to train 4 models. If we only have one GPU, we can set the `group_size` to 4. If `group_size` is set to 1, 4 models will be trained on one GPU at the same time, as there is no job scheduling system. Finally, the environment variables can be activated by `source_list`. In this example, `“source /home/user1234/deepmd.env”` is executed before `“dp”` to load the environment variables necessary to perform the training task.

### Perform model\_devi task at a local Slurm cluster

In this example, we perform the `model_devi` task at a local Slurm workstation.

```
"model_devi":
{
  "command": "lmp",
  "machine": {
    "context_type": "local",
    "batch_type": "Slurm",
    "local_root": " ./",
    "remote_root": "/home/user1234/work_path"
  },
  "resources": {
    "number_node": 1,
    "cpu_per_node": 4,
    "gpu_per_node": 1,
    "queue_name": "QueueGPU",
    "custom_flags": ["#SBATCH --mem=32G"],
    "group_size": 10,
    "source_list": ["/home/user1234/lammps.env"]
  }
}
```

The `command` for the `model_devi` task in the LAMMPS is `“lmp”`.

In the machine parameter, we specify the type of job scheduling system by changing the `batch_type` to `“Slurm”`.

In the resources parameter, we specify the name of the queue to which the task is submitted by adding `queue_name`. We can add additional lines to the calculation script via the `custom_flags`. In the `model_devi` steps, there are frequently many short tasks, so we usually pack multiple tasks (e.g. 10) into a group for submission. Other parameters are similar to that of the local workstation.

### Perform fp task in a remote PBS cluster

In this example, we perform the `fp` task at a remote PBS cluster that can be accessed via SSH.

```
"fp":
{
  "command": "mpirun -n 32 vasp_std",
  "machine": {
    "context_type": "SSHContext",
    "batch_type": "PBS",
```

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```

    "local_root": "./",
    "remote_root": "/home/user1234/work_path",
    "remote_profile": {
        "hostname": "39.xxx.xx.xx",
        "username": "user1234"
    }
},
"resources": {
    "number_node": 1,
    "cpu_per_node": 32,
    "gpu_per_node": 0,
    "queue_name": "QueueCPU",
    "group_size": 5,
    "source_list": ["/home/user1234/vasp.env"]
}
}

```

VASP code is used for fp task and mpi is used for parallel computing, so “mpirun -n 32” is added to specify the number of parallel threads.

In the machine parameter, *context\_type* is modified to “SSHContext” and *batch\_type* is modified to “PBS”. It is worth noting that *remote\_root* should be set to an accessible path on the remote PBS cluster. *remote\_profile* is added to specify the information used to connect the remote cluster, including hostname, username, port, etc.

In the resources parameter, we set *gpu\_per\_node* to 0 since it is cost-effective to use the CPU for VASP calculations.

Explicit descriptions of keys in machine.json will be given in the following section.

## 4.4 dpgen run param parameters

**Note:** One can load, modify, and export the input file by using our effective web-based tool [DP-GUI](#) online or hosted using the *command line interface* `dpgen gui`. All parameters below can be set in DP-GUI. By clicking “SAVE JSON”, one can download the input file.

### run\_jdata:

type: dict  
 argument path: run\_jdata  
 param.json file

#### type\_map:

type: list[str]  
 argument path: run\_jdata/type\_map

Atom types. Reminder: The elements in param.json, type.raw and data.lmp(when using lammps) should be in the same order.

#### mass\_map:

type: list[float] | str, optional, default: auto  
 argument path: run\_jdata/mass\_map

Standard atomic weights (default: “auto”). if one want to use isotopes, or non-standard element names, chemical symbols, or atomic number in the type\_map list, please customize the mass\_map list instead of using “auto”.

**use\_ele\_temp:**

type: int, optional, default: 0  
argument path: run\_jdata/use\_ele\_temp

Currently only support fp\_style vasp.

- 0: no electron temperature.
- 1: electron temperature as frame parameter.
- 2: electron temperature as atom parameter.

**init\_data\_prefix:**

type: str, optional  
argument path: run\_jdata/init\_data\_prefix

Prefix of initial data directories.

**init\_data\_sys:**

type: list[str]  
argument path: run\_jdata/init\_data\_sys

Paths of initial data. The path can be either a system directory containing NumPy files or an HDF5 file. You may use either absolute or relative path here. Systems will be detected recursively in the directories or the HDF5 file.

**sys\_format:**

type: str, optional, default: vasp/poscar  
argument path: run\_jdata/sys\_format

Format of sys\_configs.

**init\_batch\_size:**

type: list[typing.Union[int, str]] | str, optional  
argument path: run\_jdata/init\_batch\_size

Each number is the batch\_size of corresponding system for training in init\_data\_sys. One recommended rule for setting the sys\_batch\_size and init\_batch\_size is that batch\_size multiply number of atoms of the structure should be larger than 32. If set to auto, batch size will be 32 divided by number of atoms. This argument will not override the mixed batch size in *default\_training\_param*.

**sys\_configs\_prefix:**

type: str, optional  
argument path: run\_jdata/sys\_configs\_prefix

Prefix of sys\_configs.

**sys\_configs:**

type: list[list[str]]  
argument path: run\_jdata/sys\_configs

2D list. Containing directories of structures to be explored in iterations for each system. Wildcard characters are supported here.

**sys\_batch\_size:**

type: list[typing.Union[int, str]], optional

argument path: `run_jdata/sys_batch_size`

Each number is the `batch_size` for training of corresponding system in `sys_configs`. If set to `auto`, batch size will be 32 divided by number of atoms. This argument will not override the mixed batch size in *default\_training\_param*.

**numb\_models:**

type: `int`

argument path: `run_jdata/numb_models`

Number of models to be trained in `00.train`. 4 is recommend.

**training\_iter0\_model\_path:**

type: `list[str]`, optional

argument path: `run_jdata/training_iter0_model_path`

The model used to init the first iter training. Number of element should be equal to `numb_models`.

**training\_init\_model:**

type: `bool`, optional

argument path: `run_jdata/training_init_model`

Iteration > 0, the model parameters will be initilized from the model trained at the previous iteration. Iteration == 0, the model parameters will be initialized from `training_iter0_model_path`.

**default\_training\_param:**

type: `dict`

argument path: `run_jdata/default_training_param`

Training parameters for `deepmd-kit` in `00.train`. You can find instructions from [DeePMD-kit documentation](#).

**dp\_train\_skip\_neighbor\_stat:**

type: `bool`, optional, default: `False`

argument path: `run_jdata/dp_train_skip_neighbor_stat`

Append `-skip-neighbor-stat` flag to `dp train`.

**dp\_compress:**

type: `bool`, optional, default: `False`

argument path: `run_jdata/dp_compress`

Use `dp compress` to compress the model.

**training\_reuse\_iter:**

type: `int | NoneType`, optional

argument path: `run_jdata/training_reuse_iter`

The minimal index of iteration that continues training models from old models of last iteration.

**training\_reuse\_old\_ratio:**

type: `str | float`, optional, default: `auto`

argument path: `run_jdata/training_reuse_old_ratio`

The probability proportion of old data during training. It can be:

- `float`: directly assign the probability of old data;



- *auto:f*: automatic probability, where f is the new-to-old ratio;
- *auto*: equivalent to *auto:10*.

This option is only adopted when continuing training models from old models. This option will override default parameters.

#### **training\_reuse\_num\_steps:**

type: int | NoneType, optional, default: None, alias: *training\_reuse\_stop\_batch*  
argument path: run\_jdata/training\_reuse\_num\_steps

Number of training batch. This option is only adopted when continuing training models from old models. This option will override default parameters.

#### **training\_reuse\_start\_lr:**

type: float | NoneType, optional, default: None  
argument path: run\_jdata/training\_reuse\_start\_lr

The learning rate the start of the training. This option is only adopted when continuing training models from old models. This option will override default parameters.

#### **training\_reuse\_start\_pref\_e:**

type: int | float | NoneType, optional, default: None  
argument path: run\_jdata/training\_reuse\_start\_pref\_e

The prefactor of energy loss at the start of the training. This option is only adopted when continuing training models from old models. This option will override default parameters.

#### **training\_reuse\_start\_pref\_f:**

type: int | float | NoneType, optional, default: None  
argument path: run\_jdata/training\_reuse\_start\_pref\_f

The prefactor of force loss at the start of the training. This option is only adopted when continuing training models from old models. This option will override default parameters.

#### **model\_devi\_activation\_func:**

type: list[list[str]] | NoneType, optional  
argument path: run\_jdata/model\_devi\_activation\_func

The activation function in the model. The shape of list should be (N\_models, 2), where 2 represents the embedding and fitting network. This option will override default parameters.

#### **srtab\_file\_path:**

type: str, optional  
argument path: run\_jdata/srtab\_file\_path

The path of the table for the short-range pairwise interaction which is needed when using DP-ZBL potential

#### **one\_h5:**

type: bool, optional, default: False  
argument path: run\_jdata/one\_h5

When using DeePMD-kit, all of the input data will be merged into one HDF5 file.

**training\_init\_frozen\_model:**

type: list[str], optional

argument path: run\_jdata/training\_init\_frozen\_model

At iteration 0, initialize the model parameters from the given frozen models. Number of element should be equal to numb\_models.

**training\_finetune\_model:**

type: list[str], optional

argument path: run\_jdata/training\_finetune\_model

At iteration 0, finetune the model parameters from the given frozen models. Number of element should be equal to numb\_models.

**fp\_task\_max:**

type: int

argument path: run\_jdata/fp\_task\_max

Maximum number of structures to be calculated in each system in 02.fp of each iteration. If the number of candidate structures exceeds *fp\_task\_max*, *fp\_task\_max* structures will be randomly picked from the candidates and labeled.

**fp\_task\_min:**

type: int

argument path: run\_jdata/fp\_task\_min

Skip the training in the next iteration if the number of structures is no more than *fp\_task\_min*.

**fp\_accurate\_threshold:**

type: float, optional

argument path: run\_jdata/fp\_accurate\_threshold

If the accurate ratio is larger than this number, no fp calculation will be performed, i.e. *fp\_task\_max* = 0.

**fp\_accurate\_soft\_threshold:**

type: float, optional

argument path: run\_jdata/fp\_accurate\_soft\_threshold

If the accurate ratio is between this number and *fp\_accurate\_threshold*, the *fp\_task\_max* linearly decays to zero.

**fp\_cluster\_vacuum:**

type: float, optional

argument path: run\_jdata/fp\_cluster\_vacuum

If the vacuum size is smaller than this value, this cluster will not be chosen for labeling.

**detailed\_report\_make\_fp:**

type: bool, optional, default: True

argument path: run\_jdata/detailed\_report\_make\_fp

If set to true, a detailed report will be generated for each iteration.

**ratio\_failed:**

type: float, optional

argument path: `run_jdata/ratio_failed`

Check the ratio of unsuccessfully terminated jobs. If too many FP tasks are not converged, `RuntimeError` will be raised.

Depending on the value of `model_devi_engine`, different sub args are accepted.

#### **model\_devi\_engine:**

type: `str` (flag key), default: `lammps`

argument path: `run_jdata/model_devi_engine`

possible choices: `lammps`, `amber`, `calypso`, `gromacs`

Engine for the model deviation task.

When `model_devi_engine` is set to `lammps`:

#### **LAMMPS**

#### **model\_devi\_jobs:**

type: `list`

argument path: `run_jdata[model_devi_engine=lammps]/model_devi_jobs`

Settings for exploration in `01.model_devi`. Each dict in the list corresponds to one iteration. The index of `model_devi_jobs` exactly accord with index of iterations

This argument takes a list with each element containing the following:

#### **template:**

type: `dict`, optional

argument path:

`run_jdata[model_devi_engine=lammps]/model_devi_jobs/template`

Give an input file template for the supported engine software adopted in `01.model_devi`. Through user-defined template, any freedom (function) that is permitted by the engine software could be inherited (invoked) in the workflow.

#### **lmp:**

type: `str`, optional

argument path: `run_jdata[model_devi_engine=lammps]/model_devi_jobs/template/lmp`

The path to input.lammps template. Instructions can be found in [LAMMPS documentation](#).

#### **plm:**

type: `str`, optional

argument path: `run_jdata[model_devi_engine=lammps]/model_devi_jobs/template/plm`

The path to input.plumed template. Instructions can be found in [PLUMED documentation](#).

#### **rev\_mat:**

type: `dict`, optional

argument path: `run_jdata[model_devi_engine=lammps]/model_devi_jobs/rev_mat`

revise matrix for revising variable(s) defined in the template into the specific values (iteration-resolved). Values will be broadcasted for all tasks within the iteration invoking this key.

**lmp:**

type: dict, optional

argument path:

run\_jdata[model\_devi\_engine=lammps]/  
model\_devi\_jobs/rev\_mat/lmp

revise matrix for revising variable(s) defined in the lammps template into the specific values (iteration-resolved).

**plm:**

type: dict, optional

argument path:

run\_jdata[model\_devi\_engine=lammps]/  
model\_devi\_jobs/rev\_mat/plm

revise matrix for revising variable(s) defined in the plumed template into specific values(iteration-resolved)

**sys\_rev\_mat:**

type: dict, optional

argument path: run\_jdata[model\_devi\_engine=lammps]/  
model\_devi\_jobs/sys\_rev\_mat

system-resolved revise matrix for revising variable(s) defined in the template into specific values. Values should be individually assigned to each system adopted by this iteration, through a dictionary where first-level keys are values of sys\_idx of this iteration.

**sys\_idx:**

type: list[int]

argument path: run\_jdata[model\_devi\_engine=lammps]/  
model\_devi\_jobs/sys\_idx

Systems to be selected as the initial structure of MD and be explored. The index corresponds exactly to the sys\_configs.

**temps:**

type: list[float], optional

argument path: run\_jdata[model\_devi\_engine=lammps]/  
model\_devi\_jobs/temps

Temperature (K) in MD.

**press:**

type: list[float], optional

argument path: run\_jdata[model\_devi\_engine=lammps]/  
model\_devi\_jobs/press

Pressure (Bar) in MD. Required when ensemble is npt.

**trj\_freq:**

type: int

argument path: run\_jdata[model\_devi\_engine=lammps]/  
model\_devi\_jobs/trj\_freq

Frequency of trajectory saved in MD.

**nsteps:**

type: int, optional  
 argument path: `run_jdata[model_devi_engine=lammps]/model_devi_jobs/nsteps`  
 Running steps of MD. It is not optional when not using a template.

**nbeads:**

type: int, optional  
 argument path: `run_jdata[model_devi_engine=lammps]/model_devi_jobs/nbeads`  
 Number of beads in PIMD. If not given, classical MD will be performed. Only supported for LAMMPS version  $\geq 20230615$ .

**ensemble:**

type: str, optional  
 argument path: `run_jdata[model_devi_engine=lammps]/model_devi_jobs/ensemble`  
 Determining which ensemble used in MD, options include “npt” and “nvt”. It is not optional when not using a template.

**neidelay:**

type: int, optional  
 argument path: `run_jdata[model_devi_engine=lammps]/model_devi_jobs/neidelay`  
 delay building until this many steps since last build.

**taut:**

type: float, optional  
 argument path: `run_jdata[model_devi_engine=lammps]/model_devi_jobs/taut`  
 Coupling time of thermostat (ps).

**taup:**

type: float, optional  
 argument path: `run_jdata[model_devi_engine=lammps]/model_devi_jobs/taup`  
 Coupling time of barostat (ps).

**model\_devi\_f\_trust\_lo:**

type: dict | float, optional  
 argument path: `run_jdata[model_devi_engine=lammps]/model_devi_jobs/model_devi_f_trust_lo`  
 Lower bound of forces for the selection. If dict, should be set for each index in `sys_idx`, respectively.

**model\_devi\_f\_trust\_hi:**

type: dict | float, optional  
argument path: run\_jdata[model\_devi\_engine=lammps]/  
model\_devi\_jobs/model\_devi\_f\_trust\_hi

Upper bound of forces for the selection. If dict, should be set for each index in sys\_idx, respectively.

**model\_devi\_v\_trust\_lo:**

type: dict | float, optional  
argument path: run\_jdata[model\_devi\_engine=lammps]/  
model\_devi\_jobs/model\_devi\_v\_trust\_lo

Lower bound of virial for the selection. If dict, should be set for each index in sys\_idx, respectively. Should be used with DeePMD-kit v2.x.

**model\_devi\_v\_trust\_hi:**

type: dict | float, optional  
argument path: run\_jdata[model\_devi\_engine=lammps]/  
model\_devi\_jobs/model\_devi\_v\_trust\_hi

Upper bound of virial for the selection. If dict, should be set for each index in sys\_idx, respectively. Should be used with DeePMD-kit v2.x.

**model\_devi\_dt:**

type: float  
argument path: run\_jdata[model\_devi\_engine=lammps]/model\_devi\_dt

Timestep for MD. 0.002 is recommend.

**model\_devi\_skip:**

type: int  
argument path:  
run\_jdata[model\_devi\_engine=lammps]/model\_devi\_skip

Number of structures skipped for fp in each MD.

**model\_devi\_f\_trust\_lo:**

type: list[float] | dict | float  
argument path:  
run\_jdata[model\_devi\_engine=lammps]/model\_devi\_f\_trust\_lo

Lower bound of forces for the selection. If list or dict, should be set for each index in sys\_configs, respectively.

**model\_devi\_f\_trust\_hi:**

type: list[float] | dict | float  
argument path:  
run\_jdata[model\_devi\_engine=lammps]/model\_devi\_f\_trust\_hi

Upper bound of forces for the selection. If list or dict, should be set for each index in sys\_configs, respectively.

**model\_devi\_v\_trust\_lo:**

type: list[float] | dict | float, optional, default: 10000000000.0

argument path:

run\_jdata[model\_devi\_engine=lammps]/model\_devi\_v\_trust\_lo

Lower bound of virial for the selection. If list or dict, should be set for each index in sys\_configs, respectively. Should be used with DeePMD-kit v2.x.

**model\_devi\_v\_trust\_hi:**

type: list[float] | dict | float, optional, default: 10000000000.0

argument path:

run\_jdata[model\_devi\_engine=lammps]/model\_devi\_v\_trust\_hi

Upper bound of virial for the selection. If list or dict, should be set for each index in sys\_configs, respectively. Should be used with DeePMD-kit v2.x.

**model\_devi\_adapt\_trust\_lo:**

type: bool, optional

argument path:

run\_jdata[model\_devi\_engine=lammps]/model\_devi\_adapt\_trust\_lo

Adaptively determines the lower trust levels of force and virial. This option should be used together with model\_devi\_numb\_candi\_f, model\_devi\_numb\_candi\_v and optionally with model\_devi\_perc\_candi\_f and model\_devi\_perc\_candi\_v. dpngen will make two sets:

- 1. From the frames with force model deviation lower than model\_devi\_f\_trust\_hi, select max(model\_devi\_numb\_candi\_f, model\_devi\_perc\_candi\_f\*n\_frames) frames with largest force model deviation.
- 2. From the frames with virial model deviation lower than model\_devi\_v\_trust\_hi, select max(model\_devi\_numb\_candi\_v, model\_devi\_perc\_candi\_v\*n\_frames) frames with largest virial model deviation.

The union of the two sets is made as candidate dataset.

**model\_devi\_numb\_candi\_f:**

type: int, optional

argument path:

run\_jdata[model\_devi\_engine=lammps]/model\_devi\_numb\_candi\_f

See model\_devi\_adapt\_trust\_lo.

**model\_devi\_numb\_candi\_v:**

type: int, optional

argument path:

run\_jdata[model\_devi\_engine=lammps]/model\_devi\_numb\_candi\_v

See model\_devi\_adapt\_trust\_lo.

**model\_devi\_perc\_candi\_f:**

type: float, optional

argument path:  
run\_jdata[model\_devi\_engine=lammps]/model\_devi\_perc\_candi\_f  
See model\_devi\_adapt\_trust\_lo.

**model\_devi\_perc\_candi\_v:**

type: float, optional  
argument path:  
run\_jdata[model\_devi\_engine=lammps]/model\_devi\_perc\_candi\_v  
See model\_devi\_adapt\_trust\_lo.

**model\_devi\_f\_avg\_relative:**

type: bool, optional  
argument path:  
run\_jdata[model\_devi\_engine=lammps]/model\_devi\_f\_avg\_relative  
Normalized the force model deviations by the RMS force magnitude along the trajectory. This key should not be used with use\_relative.

**model\_devi\_clean\_traj:**

type: bool | int, optional, default: True  
argument path:  
run\_jdata[model\_devi\_engine=lammps]/model\_devi\_clean\_traj  
If type of model\_devi\_clean\_traj is bool type then it denote whether to clean traj folders in MD since they are too large. If it is Int type, then the most recent n iterations of traj folders will be retained, others will be removed.

**model\_devi\_merge\_traj:**

type: bool, optional, default: False  
argument path:  
run\_jdata[model\_devi\_engine=lammps]/model\_devi\_merge\_traj  
If model\_devi\_merge\_traj is set as True, only all.lammpstrj will be generated, instead of lots of small traj files.

**model\_devi\_nopbc:**

type: bool, optional, default: False  
argument path:  
run\_jdata[model\_devi\_engine=lammps]/model\_devi\_nopbc  
Assume open boundary condition in MD simulations.

**model\_devi\_plumed:**

type: bool, optional, default: False  
argument path:  
run\_jdata[model\_devi\_engine=lammps]/model\_devi\_plumed

**model\_devi\_plumed\_path:**

type: bool, optional, default: False  
argument path:  
run\_jdata[model\_devi\_engine=lammps]/model\_devi\_plumed\_path



**shuffle\_poscar:**

type: bool, optional, default: False

argument path: `run_jdata[model_devi_engine=lammps]/shuffle_poscar`

Shuffle atoms of each frame before running simulations. The purpose is to sample the element occupation of alloys.

**use\_relative:**

type: bool, optional, default: False

argument path: `run_jdata[model_devi_engine=lammps]/use_relative`

Calculate relative force model deviation.

**epsilon:**

type: float, optional

argument path: `run_jdata[model_devi_engine=lammps]/epsilon`

The level parameter for computing the relative force model deviation.

**use\_relative\_v:**

type: bool, optional, default: False

argument path: `run_jdata[model_devi_engine=lammps]/use_relative_v`

Calculate relative virial model deviation.

**epsilon\_v:**

type: float, optional

argument path: `run_jdata[model_devi_engine=lammps]/epsilon_v`

The level parameter for computing the relative virial model deviation.

When `model_devi_engine` is set to `amber`:

Amber DPRc engine. The command argument in the machine file should be path to `sander`.

**model\_devi\_jobs:**

type: list

argument path: `run_jdata[model_devi_engine=amber]/model_devi_jobs`

List of dicts. The list including the dict for information of each cycle.

This argument takes a list with each element containing the following:

**sys\_idx:**

type: list[int]

argument path: `run_jdata[model_devi_engine=amber]/model_devi_jobs/sys_idx`

List of ints. List of systems to run.

**trj\_freq:**

type: int

argument path: `run_jdata[model_devi_engine=amber]/model_devi_jobs/trj_freq`

Frequency to dump trajectory.

**restart\_from\_iter:**

type: int, optional

argument path: run\_jdata[model\_devi\_engine=amber]/  
model\_devi\_jobs/restart\_from\_iter

The iteration index to restart the simulation from. If not given, the simulation is restarted from *sys\_configs*.

**low\_level:**

type: str

argument path: run\_jdata[model\_devi\_engine=amber]/low\_level

Low level method. The value will be filled into mdin file as @qm\_theory@.

**cutoff:**

type: float

argument path: run\_jdata[model\_devi\_engine=amber]/cutoff

Cutoff radius for the DPRc model.

**parm7\_prefix:**

type: str, optional

argument path: run\_jdata[model\_devi\_engine=amber]/parm7\_prefix

The path prefix to AMBER PARM7 files.

**parm7:**

type: list[str]

argument path: run\_jdata[model\_devi\_engine=amber]/parm7

List of paths to AMBER PARM7 files. Each file maps to a system.

**mdin\_prefix:**

type: str, optional

argument path: run\_jdata[model\_devi\_engine=amber]/mdin\_prefix

The path prefix to AMBER mdin template files.

**mdin:**

type: list[str]

argument path: run\_jdata[model\_devi\_engine=amber]/mdin

List of paths to AMBER mdin template files. Each files maps to a system. In the template, the following keywords will be replaced by the actual value: @freq@: freq to dump trajectory; @nstlim@: total time step to run; @qm\_region@: AMBER mask of the QM region; @qm\_theory@: The low level QM theory, such as DFTB2; @qm\_charge@: The total charge of the QM theory, such as -2; @rcut@: cutoff radius of the DPRc model; @GRAPH\_FILE0@, @GRAPH\_FILE1@, ... : graph files.

**qm\_region:**

type: list[str]

argument path: run\_jdata[model\_devi\_engine=amber]/qm\_region

List of strings. AMBER mask of the QM region. Each mask maps to a system.

**qm\_charge:**

type: `list[int]`  
 argument path: `run_jdata[model_devi_engine=amber]/qm_charge`  
 List of ints. Charge of the QM region. Each charge maps to a system.

**nsteps:**

type: `list[int]`  
 argument path: `run_jdata[model_devi_engine=amber]/nsteps`  
 List of ints. The number of steps to run. Each number maps to a system.

**r:**

type: `list[list[typing.Union[float, list[float]]]]`  
 argument path: `run_jdata[model_devi_engine=amber]/r`  
 2D or 3D list of floats. Constrict values for the enhanced sampling. The first dimension maps to systems. The second dimension maps to confs in each system. The third dimension is the constrict value. It can be a single float for 1D or list of floats for nD.

**disang\_prefix:**

type: `str`, optional  
 argument path: `run_jdata[model_devi_engine=amber]/disang_prefix`  
 The path prefix to disang prefix.

**disang:**

type: `list[str]`  
 argument path: `run_jdata[model_devi_engine=amber]/disang`  
 List of paths to AMBER disang files. Each file maps to a system. The keyword RVAL will be replaced by the constrict values, or RVAL1, RVAL2, ... for an nD system.

**model\_devi\_f\_trust\_lo:**

type: `list[float] | dict | float`  
 argument path:  
`run_jdata[model_devi_engine=amber]/model_devi_f_trust_lo`  
 Lower bound of forces for the selection. If dict, should be set for each index in `sys_idx`, respectively.

**model\_devi\_f\_trust\_hi:**

type: `list[float] | dict | float`  
 argument path:  
`run_jdata[model_devi_engine=amber]/model_devi_f_trust_hi`  
 Upper bound of forces for the selection. If dict, should be set for each index in `sys_idx`, respectively.

When `model_devi_engine` is set to `calypso`:

TODO: add doc

When `model_devi_engine` is set to `gromacs`:

TODO: add doc

Depending on the value of *fp\_style*, different sub args are accepted.

**fp\_style:**

type: str (flag key)  
argument path: run\_jdata/fp\_style  
possible choices: *vasp*, *gaussian*, *siesta*, *cp2k*, *abacus*, *amber/diff*,  
*pwmat*, *pwscf*, *custom*  
Software for First Principles.

When *fp\_style* is set to *vasp*:

**fp\_pp\_path:**

type: str  
argument path: run\_jdata[fp\_style=vasp]/fp\_pp\_path  
Directory of psuedo-potential file to be used for 02.fp exists.

**fp\_pp\_files:**

type: list[str]  
argument path: run\_jdata[fp\_style=vasp]/fp\_pp\_files  
Psuedo-potential file to be used for 02.fp. Note that the order of elements should correspond to the order in type\_map.

**fp\_incar:**

type: str  
argument path: run\_jdata[fp\_style=vasp]/fp\_incar  
Input file for VASP. INCAR must specify KSPACING and KGAMMA.

**fp\_aniso\_kspacing:**

type: list[float], optional  
argument path: run\_jdata[fp\_style=vasp]/fp\_aniso\_kspacing  
Set anisotropic kspacing. Usually useful for 1-D or 2-D materials. Only support VASP. If it is setting the KSPACING key in INCAR will be ignored.

**cvasp:**

type: bool, optional  
argument path: run\_jdata[fp\_style=vasp]/cvasp  
If cvasp is true, DP-GEN will use Custodian to help control VASP calculation.

**fp\_skip\_bad\_box:**

type: str, optional  
argument path: run\_jdata[fp\_style=vasp]/fp\_skip\_bad\_box  
Skip the configurations that are obviously unreasonable before 02.fp

When *fp\_style* is set to *gaussian*:

**use\_clusters:**

type: bool, optional, default: False

argument path: `run_jdata[fp_style=gaussian]/use_clusters`

If set to true, clusters will be taken instead of the whole system.

**cluster\_cutoff:**

type: float, optional

argument path: `run_jdata[fp_style=gaussian]/cluster_cutoff`

The soft cutoff radius of clusters if *use\_clusters* is set to true. Molecules will be taken as whole even if part of atoms is out of the cluster. Use *cluster\_cutoff\_hard* to only take atoms within the hard cutoff radius.

**cluster\_cutoff\_hard:**

type: float, optional

argument path: `run_jdata[fp_style=gaussian]/cluster_cutoff_hard`

The hard cutoff radius of clusters if *use\_clusters* is set to true. Outside the hard cutoff radius, atoms will not be taken even if they are in a molecule where some atoms are within the cutoff radius.

**cluster\_minify:**

type: bool, optional, default: False

argument path: `run_jdata[fp_style=gaussian]/cluster_minify`

If enabled, when an atom within the soft cutoff radius connects a single bond with a non-hydrogen atom out of the soft cutoff radius, the outer atom will be replaced by a hydrogen atom. When the outer atom is a hydrogen atom, the outer atom will be kept. In this case, other atoms out of the soft cutoff radius will be removed.

**fp\_params:**

type: dict

argument path: `run_jdata[fp_style=gaussian]/fp_params`

Parameters for Gaussian calculation.

**keywords:**

type: list[str] | str

argument path:

`run_jdata[fp_style=gaussian]/fp_params/keywords`

Keywords for Gaussian input, e.g. `force b3lyp/6-31g**`. If a list, run multiple steps.

**multiplicity:**

type: str | int, optional, default: auto

argument path:

`run_jdata[fp_style=gaussian]/fp_params/multiplicity`

Spin multiplicity for Gaussian input. If *auto*, multiplicity will be detected automatically, with the following rules: when *fragment\_guesses=True*, multiplicity will +1 for each radical, and +2 for each oxygen molecule; when *fragment\_guesses=False*, multiplicity will be 1 or 2, but +2 for each oxygen molecule.

**nproc:**

type: int  
argument path:  
run\_jdata[fp\_style=gaussian]/fp\_params/nproc

The number of processors for Gaussian input.

**charge:**

type: int, optional, default: 0  
argument path:  
run\_jdata[fp\_style=gaussian]/fp\_params/charge

Molecule charge. Only used when charge is not provided by the system.

**fragment\_guesses:**

type: bool, optional, default: False  
argument path: run\_jdata[fp\_style=gaussian]/fp\_params/  
fragment\_guesses

Initial guess generated from fragment guesses. If True, *multiplicity* should be *auto*.

**basis\_set:**

type: str, optional  
argument path:  
run\_jdata[fp\_style=gaussian]/fp\_params/basis\_set

Custom basis set.

**keywords\_high\_multiplicity:**

type: str, optional  
argument path: run\_jdata[fp\_style=gaussian]/fp\_params/  
keywords\_high\_multiplicity

Keywords for points with multiple raicals. *multiplicity* should be *auto*. If not set, fallback to normal keywords.

When `fp_style` is set to `siesta`:

**use\_clusters:**

type: bool, optional  
argument path: run\_jdata[fp\_style=siesta]/use\_clusters

If set to true, clusters will be taken instead of the whole system. This option does not work with DeePMD-kit 0.x.

**cluster\_cutoff:**

type: float, optional  
argument path: run\_jdata[fp\_style=siesta]/cluster\_cutoff

The cutoff radius of clusters if `use_clusters` is set to true.

**fp\_params:**

type: dict

argument path: `run_jdata[fp_style=siesta]/fp_params`

Parameters for siesta calculation.

**ecut:**

type: int

argument path:

`run_jdata[fp_style=siesta]/fp_params/ecut`

Define the plane wave cutoff for grid.

**ediff:**

type: float

argument path:

`run_jdata[fp_style=siesta]/fp_params/ediff`

Tolerance of Density Matrix.

**kspacing:**

type: float

argument path:

`run_jdata[fp_style=siesta]/fp_params/kspacing`

Sample factor in Brillouin zones.

**mixingWeight:**

type: float

argument path:

`run_jdata[fp_style=siesta]/fp_params/mixingWeight`

Proportion a of output Density Matrix to be used for the input Density Matrix of next SCF cycle (linear mixing).

**NumberPulay:**

type: int

argument path:

`run_jdata[fp_style=siesta]/fp_params/NumberPulay`

Controls the Pulay convergence accelerator.

**fp\_pp\_path:**

type: str

argument path: `run_jdata[fp_style=siesta]/fp_pp_path`

Directory of psuedo-potential or numerical orbital files to be used for 02.fp exists.

**fp\_pp\_files:**

type: list[str]

argument path: `run_jdata[fp_style=siesta]/fp_pp_files`

Psuedo-potential file to be used for 02.fp. Note that the order of elements should correspond to the order in type\_map.

When `fp_style` is set to `cp2k`:

**user\_fp\_params:**

type: dict, optional, alias: `fp_params`

argument path: `run_jdata[fp_style=cp2k]/user_fp_params`

Parameters for cp2k calculation. find detail in [manual.cp2k.org](http://manual.cp2k.org). only the kind section must be set before use. we assume that you have basic knowledge for cp2k input.

**external\_input\_path:**

type: str, optional

argument path: `run_jdata[fp_style=cp2k]/external_input_path`

Conflict with key: `user_fp_params`. enable the template input provided by user. some rules should be followed, read the following text in detail:

1. One must present a KEYWORD ABC in the section CELL so that the DP-GEN can replace the cell on-the-fly.
2. One need to add these lines under FORCE\_EVAL section to print forces and stresses:

```
STRESS_TENSOR ANALYTICAL
&PRINT
  &FORCES ON
  &END FORCES
  &STRESS_TENSOR ON
  &END STRESS_TENSOR
&END PRINT
```

When `fp_style` is set to `abacus`:

**fp\_pp\_path:**

type: str

argument path: `run_jdata[fp_style=abacus]/fp_pp_path`

Directory of psuedo-potential or numerical orbital files to be used for 02.fp exists.

**fp\_pp\_files:**

type: list[str]

argument path: `run_jdata[fp_style=abacus]/fp_pp_files`

Psuedo-potential file to be used for 02.fp. Note that the order of elements should correspond to the order in `type_map`.

**fp\_orb\_files:**

type: list[str], optional

argument path: `run_jdata[fp_style=abacus]/fp_orb_files`

numerical orbital file to be used for 02.fp when using LCAO basis. Note that the order of elements should correspond to the order in `type_map`.

**fp\_incar:**

type: str, optional



argument path: `run_jdata[fp_style=abacus]/fp_incar`

Input file for ABACUS. This is optional but the priority is lower than `user_fp_params`, and you should not set `user_fp_params` if you want to use `fp_incar`.

**fp\_kpt\_file:**

type: str, optional

argument path: `run_jdata[fp_style=abacus]/fp_kpt_file`

KPT file for ABACUS. If the “kspacing” or “gamma\_only=1” is defined in INPUT or “k\_points” is defined, `fp_kpt_file` will be ignored.

**fp\_dpks\_descriptor:**

type: str, optional

argument path: `run_jdata[fp_style=abacus]/fp_dpks_descriptor`

DeePKS descriptor file name. The file should be in pseudopotential directory.

**user\_fp\_params:**

type: dict, optional

argument path: `run_jdata[fp_style=abacus]/user_fp_params`

Set the key and value of INPUT.

**k\_points:**

type: list[int], optional

argument path: `run_jdata[fp_style=abacus]/k_points`

Monkhorst-Pack k-grids setting for generating KPT file of ABACUS, such as: [1,1,1,0,0,0]. NB: if “kspacing” or “gamma\_only=1” is defined in INPUT, `k_points` will be ignored.

When `fp_style` is set to `amber/diff`:

Amber/diff style for DPRc models. Note: this fp style only supports to be used with `model_devi_engine` *amber*, where some arguments are reused. The command argument in the machine file should be path to sander. One should also install `dpamber` and make it visible in the PATH.

**high\_level:**

type: str

argument path: `run_jdata[fp_style=amber/diff]/high_level`

High level method. The value will be filled into `mdin` template as `@qm_theory@`.

**fp\_params:**

type: dict

argument path: `run_jdata[fp_style=amber/diff]/fp_params`

Parameters for FP calculation.

**high\_level\_mdin:**

type: str

argument path: `run_jdata[fp_style=amber/diff]/fp_params/high_level_mdin`

Path to high-level AMBER mdin template file. `%qm_theory%`, `%qm_region%`, and `%qm_charge%` will be replaced.

**low\_level\_mdin:**

type: str

argument path: `run_jdata[fp_style=amber/diff]/fp_params/low_level_mdin`

Path to low-level AMBER mdin template file. `%qm_theory%`, `%qm_region%`, and `%qm_charge%` will be replaced.

When `fp_style` is set to `pwmat`:

TODO: add doc

When `fp_style` is set to `pwscf`:

**fp\_pp\_path:**

type: str

argument path: `run_jdata[fp_style=pwscf]/fp_pp_path`

Directory of psuedo-potential file to be used for 02.fp exists.

**fp\_pp\_files:**

type: list[str]

argument path: `run_jdata[fp_style=pwscf]/fp_pp_files`

Psuedo-potential file to be used for 02.fp. Note that the order of elements should correspond to the order in `type_map`.

**fp\_params:**

type: dict, optional

argument path: `run_jdata[fp_style=pwscf]/fp_params`

Parameters for pwscf calculation. It has lower priority than `user_fp_params`.

**ecut:**

type: float

argument path:

`run_jdata[fp_style=pwscf]/fp_params/ecut`

`ecutwfc` in pwscf.

**ediff:**

type: float

argument path:

`run_jdata[fp_style=pwscf]/fp_params/ediff`

`conv_thr` and `ts_vdw_econv_thr` in pwscf.

**smearing:**

type: str

argument path:  
`run_jdata[fp_style=pwscf]/fp_params/smearing`  
 smearing in pwscf.

**sigma:**

type: float  
 argument path:  
`run_jdata[fp_style=pwscf]/fp_params/sigma`  
 degauss in pwscf.

**kspacing:**

type: float  
 argument path:  
`run_jdata[fp_style=pwscf]/fp_params/kspacing`  
 The spacing between kpoints. Helps to determin KPOINTS in pwscf.

**user\_fp\_params:**

type: dict, optional  
 argument path: `run_jdata[fp_style=pwscf]/user_fp_params`  
 Parameters for pwscf calculation. Find details at [https://www.quantum-espresso.org/Doc/INPUT\\_PW.html](https://www.quantum-espresso.org/Doc/INPUT_PW.html). When `user_fp_params` is set, the settings in `fp_params` will be ignored. If one wants to use `user_fp_params`, `kspacing` must be set in `user_fp_params`. `kspacing` is the spacing between kpoints, and helps to determin KPOINTS in pwscf.

When `fp_style` is set to custom:

Custom FP code. You need to provide the input and output file format and name. The command argument in the machine file should be the script to run custom FP codes. The extra forward and backward files can be defined in the machine file.

**fp\_params:**

type: dict  
 argument path: `run_jdata[fp_style=custom]/fp_params`  
 Parameters for FP calculation.

**input\_fmt:**

type: str  
 argument path:  
`run_jdata[fp_style=custom]/fp_params/input_fmt`  
 Input dpdata format of the custom FP code. Such format should only need the first argument as the file name.

**input\_fn:**

type: str  
 argument path:  
`run_jdata[fp_style=custom]/fp_params/input_fn`  
 Input file name of the custom FP code.

**output\_fmt:**

type: str  
argument path:  
run\_jdata[fp\_style=custom]/fp\_params/output\_fmt  
Output dpdata format of the custom FP code. Such format should only need the first argument as the file name.

**output\_fn:**

type: str  
argument path:  
run\_jdata[fp\_style=custom]/fp\_params/output\_fn  
Output file name of the custom FP code.

## 4.5 dpgen run machine parameters

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**Note:** One can load, modify, and export the input file by using our effective web-based tool [DP-GUI](#) online or hosted using the *command line interface* `dpgen gui`. All parameters below can be set in DP-GUI. By clicking “SAVE JSON”, one can download the input file.

---

**run\_mdata:**

type: dict  
argument path: run\_mdata  
machine.json file

**api\_version:**

type: str, optional, default: 1.0  
argument path: run\_mdata/api\_version  
Please set to 1.0

**deepmd\_version:**

type: str, optional, default: 2  
argument path: run\_mdata/deepmd\_version  
DeePMD-kit version, e.g. 2.1.3

**train:**

type: dict  
argument path: run\_mdata/train  
Parameters of command, machine, and resources for train

**command:**

type: str  
argument path: run\_mdata/train/command  
Command of a program.

**machine:**

type: dict

argument path: run\_mdata/train/machine

**batch\_type:**

type: str

argument path:

run\_mdata/train/machine/batch\_type

The batch job system type. Option: Fugaku, Slurm, DistributedShell, SGE, PBS, Torque, SlurmJobArray, Shell, Bohrium, LSF, OpenAPI

**local\_root:**

type: str | NoneType

argument path:

run\_mdata/train/machine/local\_root

The dir where the tasks and relating files locate. Typically the project dir.

**remote\_root:**

type: str | NoneType, optional

argument path:

run\_mdata/train/machine/remote\_root

The dir where the tasks are executed on the remote machine. Only needed when context is not lazy-local.

**clean\_asynchronously:**

type: bool, optional, default: False

argument path: run\_mdata/train/machine/

clean\_asynchronously

Clean the remote directory asynchronously after the job finishes.

Depending on the value of *context\_type*, different sub args are accepted.

**context\_type:**

type: str (flag key)

argument path:

run\_mdata/train/machine/context\_type

possible choices: *BohriumContext*, *LocalContext*, *LazyLocalContext*, *SSHContext*, *HDFSContext*, *OpenAPIContext*

The connection used to remote machine. Option: OpenAPIContext, SSHContext, BohriumContext, HDFSContext, LazyLocalContext, LocalContext

When *context\_type* is set to *BohriumContext* (or its aliases *bohriumcontext*, *Bohrium*, *bohrium*, *DpCloudServerContext*, *dpcloudservercontext*, *DpCloudServer*, *dpcloudserver*, *LebesgueContext*, *lebesguecontext*, *Lebesgue*, *lebesgue*):

**remote\_profile:**

type: dict

argument path: run\_mdata/train/

machine[BohriumContext]/remote\_profile

The information used to maintain the connection with remote machine.

**email:**

type: str, optional

argument path: run\_mdata/train/

machine[BohriumContext]/

remote\_profile/email

Email

**password:**

type: str, optional

argument path: run\_mdata/train/

machine[BohriumContext]/

remote\_profile/password

Password

**program\_id:**

type: int, alias: *project\_id*

argument path: run\_mdata/train/

machine[BohriumContext]/

remote\_profile/program\_id

Program ID

**retry\_count:**

type: NoneType | int, optional, default: 2

argument path: run\_mdata/train/

machine[BohriumContext]/

remote\_profile/retry\_count

The retry count when a job is terminated

**ignore\_exit\_code:**

type: bool, optional, default: True

argument path: run\_mdata/train/

machine[BohriumContext]/

remote\_profile/ignore\_exit\_code

**The job state will be marked as finished if the exit code is non-zero when set to True.**

**Otherwise,**

the job state will be designated as terminated.

**keep\_backup:**

type: bool, optional

argument path: run\_mdata/train/

machine[BohriumContext]/

remote\_profile/keep\_backup

keep download and upload zip

**input\_data:**

type: dict  
 argument path: run\_mdata/train/  
 machine[BohriumContext]/  
 remote\_profile/input\_data

Configuration of job

When `context_type` is set to `LocalContext` (or its aliases `localcontext`, `Local`, `local`):

**remote\_profile:**

type: dict, optional  
 argument path: run\_mdata/train/  
 machine[LocalContext]/remote\_profile

The information used to maintain the connection with remote machine. This field is empty for this context.

When `context_type` is set to `LazyLocalContext` (or its aliases `lazylocalcontext`, `LazyLocal`, `lazylocal`):

**remote\_profile:**

type: dict, optional  
 argument path: run\_mdata/train/  
 machine[LazyLocalContext]/  
 remote\_profile

The information used to maintain the connection with remote machine. This field is empty for this context.

When `context_type` is set to `SSHContext` (or its aliases `sshcontext`, `SSH`, `ssh`):

**remote\_profile:**

type: dict  
 argument path: run\_mdata/train/  
 machine[SSHContext]/remote\_profile

The information used to maintain the connection with remote machine.

**hostname:**

type: str  
 argument path: run\_mdata/train/  
 machine[SSHContext]/  
 remote\_profile/hostname

hostname or ip of ssh connection.

**username:**

type: str

argument path: run\_mdata/train/  
machine[SSHContext]/  
remote\_profile/username

username of target linux system

**password:**

type: str, optional  
argument path: run\_mdata/train/  
machine[SSHContext]/  
remote\_profile/password  
  
(deprecated) password of linux system.  
Please use [SSH keys](#) instead to improve  
security.

**port:**

type: int, optional, default: 22  
argument path: run\_mdata/train/  
machine[SSHContext]/  
remote\_profile/port  
  
ssh connection port.

**key\_filename:**

type: str | NoneType, optional, default:  
None  
argument path: run\_mdata/train/  
machine[SSHContext]/  
remote\_profile/key\_filename  
  
key filename used by ssh connection. If  
left None, find key in ~/.ssh or use pass-  
word for login

**passphrase:**

type: str | NoneType, optional, default:  
None  
argument path: run\_mdata/train/  
machine[SSHContext]/  
remote\_profile/passphrase  
  
passphrase of key used by ssh connection

**timeout:**

type: int, optional, default: 10  
argument path: run\_mdata/train/  
machine[SSHContext]/  
remote\_profile/timeout  
  
timeout of ssh connection

**totp\_secret:**

type: str | NoneType, optional, default:  
None  
argument path: run\_mdata/train/  
machine[SSHContext]/  
remote\_profile/totp\_secret



Time-based one time password secret. It should be a base32-encoded string extracted from the 2D code.

**tar\_compress:**

type: bool, optional, default: True  
 argument path: run\_mdata/train/  
 machine[SSHContext]/  
 remote\_profile/tar\_compress

The archive will be compressed in upload and download if it is True. If not, compression will be skipped.

**look\_for\_keys:**

type: bool, optional, default: True  
 argument path: run\_mdata/train/  
 machine[SSHContext]/  
 remote\_profile/look\_for\_keys

enable searching for discoverable private key files in ~/.ssh/

When `context_type` is set to `HDFSContext` (or its aliases `hdfscontext`, `HDFS`, `hdfs`):

**remote\_profile:**

type: dict, optional  
 argument path: run\_mdata/train/  
 machine[HDFSContext]/remote\_profile

The information used to maintain the connection with remote machine. This field is empty for this context.

When `context_type` is set to `OpenAPIContext` (or its aliases `openapicontext`, `OpenAPI`, `openapi`):

**remote\_profile:**

type: dict, optional  
 argument path: run\_mdata/train/  
 machine[OpenAPIContext]/remote\_profile

The information used to maintain the connection with remote machine. This field is empty for this context.

**resources:**

type: dict  
 argument path: run\_mdata/train/resources

**number\_node:**

type: int, optional, default: 1  
 argument path: run\_mdata/train/  
 resources/number\_node

The number of node need for each *job*

**cpu\_per\_node:**

type: int, optional, default: 1

argument path: run\_mdata/train/  
resources/cpu\_per\_node

cpu numbers of each node assigned to each job.

**gpu\_per\_node:**

type: int, optional, default: 0

argument path: run\_mdata/train/  
resources/gpu\_per\_node

gpu numbers of each node assigned to each job.

**queue\_name:**

type: str, optional, default: (empty string)

argument path: run\_mdata/train/  
resources/queue\_name

The queue name of batch job scheduler system.

**group\_size:**

type: int

argument path: run\_mdata/train/  
resources/group\_size

The number of *tasks* in a *job*. 0 means infinity.

**custom\_flags:**

type: typing.List[str], optional

argument path: run\_mdata/train/  
resources/custom\_flags

The extra lines pass to job submitting script  
header

**strategy:**

type: dict, optional

argument path:  
run\_mdata/train/resources/strategy

strategies we use to generation job submitting  
scripts.

**if\_cuda\_multi\_devices:**

type: bool, optional, default: False

argument path: run\_mdata/train/  
resources/strategy/  
if\_cuda\_multi\_devices

If there are multiple nvidia GPUS  
on the node, and we want to as-  
sign the tasks to different GPUS.If

true, dpdispatcher will manually export environment variable `CUDA_VISIBLE_DEVICES` to different task. Usually, this option will be used with `Task.task_need_resources` variable simultaneously.

**ratio\_unfinished:**

type: float, optional, default: 0.0

argument path:

`run_mdata/train/resources/  
strategy/ratio_unfinished`

The ratio of *tasks* that can be unfinished.

**customized\_script\_header\_template\_file:**

type: str, optional

argument path: `run_mdata/train/  
resources/strategy/`

`customized_script_header_template_file`

The customized template file to generate job submitting script header, which overrides the default file.

**para\_deg:**

type: int, optional, default: 1

argument path:

`run_mdata/train/resources/para_deg`

Decide how many tasks will be run in parallel.

**source\_list:**

type: `typing.List[str]`, optional, default: []

argument path: `run_mdata/train/  
resources/source_list`

The env file to be sourced before the command execution.

**module\_purge:**

type: bool, optional, default: False

argument path: `run_mdata/train/  
resources/module_purge`

Remove all modules on HPC system before module load (`module_list`)

**module\_unload\_list:**

type: `typing.List[str]`, optional, default: []

argument path: `run_mdata/train/  
resources/module_unload_list`

The modules to be unloaded on HPC system before submitting jobs

**module\_list:**

type: `typing.List[str]`, optional, default: `[]`

argument path: `run_mdata/train/resources/module_list`

The modules to be loaded on HPC system before submitting jobs

**envs:**

type: `dict`, optional, default: `{}`

argument path: `run_mdata/train/resources/envs`

The environment variables to be exported on before submitting jobs

**prepend\_script:**

type: `typing.List[str]`, optional, default: `[]`

argument path: `run_mdata/train/resources/prepend_script`

Optional script run before jobs submitted.

**append\_script:**

type: `typing.List[str]`, optional, default: `[]`

argument path: `run_mdata/train/resources/append_script`

Optional script run after jobs submitted.

**wait\_time:**

type: `float | int`, optional, default: `0`

argument path: `run_mdata/train/resources/wait_time`

The waiting time in second after a single *task* submitted

Depending on the value of *batch\_type*, different sub args are accepted.

**batch\_type:**

type: `str` (flag key)

argument path: `run_mdata/train/resources/batch_type`

possible choices: *Bohrium*, *Fugaku*, *Slurm*, *SlurmJobArray*, *Shell*, *LSF*, *PBS*, *SGE*, *Torque*, *OpenAPI*, *DistributedShell*

The batch job system type loaded from machine/batch\_type.

When `batch_type` is set to Bohrium (or its aliases bohrium, Lebesgue, lebesgue, DpCloudServer, dpcloudserver):

**kwargs:**

type: dict, optional  
 argument path: run\_mdata/train/  
 resources[Bohrium]/kwargs

This field is empty for this batch.

When `batch_type` is set to Fugaku (or its alias fugaku):

**kwargs:**

type: dict, optional  
 argument path: run\_mdata/train/  
 resources[Fugaku]/kwargs

This field is empty for this batch.

When `batch_type` is set to Slurm (or its alias slurm):

**kwargs:**

type: dict, optional  
 argument path: run\_mdata/train/  
 resources[Slurm]/kwargs

Extra arguments.

**custom\_gpu\_line:**

type: str | NoneType, optional,  
 default: None  
 argument path: run\_mdata/train/  
 resources[Slurm]/kwargs/  
 custom\_gpu\_line  
 Custom GPU configuration, starting  
 with #SBATCH

When `batch_type` is set to SlurmJobArray (or its alias slurmjobarray):

**kwargs:**

type: dict, optional  
 argument path: run\_mdata/train/  
 resources[SlurmJobArray]/kwargs

Extra arguments.

**custom\_gpu\_line:**

type: str | NoneType, optional,  
 default: None  
 argument path: run\_mdata/train/  
 resources[SlurmJobArray]/  
 kwargs/custom\_gpu\_line

Custom GPU configuration, starting with #SBATCH

**slurm\_job\_size:**

type: int, optional, default: 1  
argument path: run\_mdata/train/  
resources[SlurmJobArray]/  
kwargs/slurm\_job\_size

Number of tasks in a Slurm job

When `batch_type` is set to Shell (or its alias `shell`):

**kwargs:**

type: dict, optional  
argument path: run\_mdata/train/  
resources[Shell]/kwargs

This field is empty for this batch.

When `batch_type` is set to LSF (or its alias `lsf`):

**kwargs:**

type: dict  
argument path: run\_mdata/train/  
resources[LSF]/kwargs

Extra arguments.

**gpu\_usage:**

type: bool, optional, default: False  
argument path: run\_mdata/train/  
resources[LSF]/kwargs/  
gpu\_usage

Choosing if GPU is used in the calculation step.

**gpu\_new\_syntax:**

type: bool, optional, default: False  
argument path: run\_mdata/train/  
resources[LSF]/kwargs/  
gpu\_new\_syntax

For LFS  $\geq$  10.1.0.3, new option -gpu for #BSUB could be used. If False, and old syntax would be used.

**gpu\_exclusive:**

type: bool, optional, default: True  
argument path: run\_mdata/train/  
resources[LSF]/kwargs/  
gpu\_exclusive

Only take effect when new syntax enabled. Control whether submit tasks in exclusive way for GPU.

**custom\_gpu\_line:**

type: str | NoneType, optional,  
 default: None  
 argument path: run\_mdata/train/  
 resources[LSF]/kwargs/  
 custom\_gpu\_line  
 Custom GPU configuration, starting  
 with #BSUB

When `batch_type` is set to PBS (or its alias `pbs`):

**kwargs:**

type: dict, optional  
 argument path: run\_mdata/train/  
 resources[PBS]/kwargs  
 This field is empty for this batch.

When `batch_type` is set to SGE (or its alias `sge`):

**kwargs:**

type: dict, optional  
 argument path: run\_mdata/train/  
 resources[SGE]/kwargs  
 This field is empty for this batch.

When `batch_type` is set to Torque (or its alias `torque`):

**kwargs:**

type: dict, optional  
 argument path: run\_mdata/train/  
 resources[Torque]/kwargs  
 This field is empty for this batch.

When `batch_type` is set to OpenAPI (or its alias `openapi`):

**kwargs:**

type: dict, optional  
 argument path: run\_mdata/train/  
 resources[OpenAPI]/kwargs  
 This field is empty for this batch.

When `batch_type` is set to DistributedShell (or its alias `distributedshell`):

**kwargs:**

type: dict, optional  
 argument path: run\_mdata/train/  
 resources[DistributedShell]/kwargs  
 This field is empty for this batch.

**user\_forward\_files:**

type: list, optional  
argument path: run\_mdata/train/user\_forward\_files  
Files to be forwarded to the remote machine.

**user\_backward\_files:**

type: list, optional  
argument path:  
run\_mdata/train/user\_backward\_files  
Files to be backwarded from the remote machine.

**model\_devi:**

type: dict  
argument path: run\_mdata/model\_devi  
Parameters of command, machine, and resources for model\_devi

**command:**

type: str  
argument path: run\_mdata/model\_devi/command  
Command of a program.

**machine:**

type: dict  
argument path: run\_mdata/model\_devi/machine

**batch\_type:**

type: str  
argument path: run\_mdata/model\_devi/  
machine/batch\_type  
The batch job system type. Option: Fugaku,  
Slurm, DistributedShell, SGE, PBS, Torque,  
SlurmJobArray, Shell, Bohrium, LSF, Ope-  
nAPI

**local\_root:**

type: str | NoneType  
argument path: run\_mdata/model\_devi/  
machine/local\_root  
The dir where the tasks and relating files locate.  
Typically the project dir.

**remote\_root:**

type: str | NoneType, optional  
argument path: run\_mdata/model\_devi/  
machine/remote\_root  
The dir where the tasks are executed on the re-  
mote machine. Only needed when context is  
not lazy-local.



**clean\_asynchronously:**

type: bool, optional, default: False  
 argument path: run\_mdata/model\_devi/  
 machine/clean\_asynchronously

Clean the remote directory asynchronously after the job finishes.

Depending on the value of *context\_type*, different sub args are accepted.

**context\_type:**

type: str (flag key)  
 argument path: run\_mdata/model\_devi/  
 machine/context\_type  
 possible choices: *BohriumContext*,  
*LocalContext*, *LazyLocalContext*,  
*SSHContext*, *HDFSContext*,  
*OpenAPIContext*

The connection used to remote machine. Option: OpenAPIContext, SSHContext, BohriumContext, HDFSContext, LazyLocalContext, LocalContext

When *context\_type* is set to *BohriumContext* (or its aliases *bohriumcontext*, *Bohrium*, *bohrium*, *DpCloudServerContext*, *dpcloudservercontext*, *DpCloudServer*, *dpcloudserver*, *LebesgueContext*, *lebesguecontext*, *Lebesgue*, *lebesgue*):

**remote\_profile:**

type: dict  
 argument path: run\_mdata/model\_devi/  
 machine[BohriumContext]/  
 remote\_profile

The information used to maintain the connection with remote machine.

**email:**

type: str, optional  
 argument path:  
 run\_mdata/model\_devi/  
 machine[BohriumContext]/  
 remote\_profile/email

Email

**password:**

type: str, optional  
 argument path:  
 run\_mdata/model\_devi/  
 machine[BohriumContext]/  
 remote\_profile/password

Password

**program\_id:**

type: int, alias: *project\_id*  
argument path:  
run\_mdata/model\_devi/  
machine[BohriumContext]/  
remote\_profile/program\_id

Program ID

**retry\_count:**

type: NoneType | int, optional,  
default: 2  
argument path:  
run\_mdata/model\_devi/  
machine[BohriumContext]/  
remote\_profile/retry\_count

The retry count when a job is terminated

**ignore\_exit\_code:**

type: bool, optional, default: True  
argument path:  
run\_mdata/model\_devi/  
machine[BohriumContext]/  
remote\_profile/  
ignore\_exit\_code

**The job state will be marked as finished if the exit code is non-zero when set to True. Otherwise,**  
the job state will be designated as terminated.

**keep\_backup:**

type: bool, optional  
argument path:  
run\_mdata/model\_devi/  
machine[BohriumContext]/  
remote\_profile/keep\_backup

keep download and upload zip

**input\_data:**

type: dict  
argument path:  
run\_mdata/model\_devi/  
machine[BohriumContext]/  
remote\_profile/input\_data

Configuration of job

When `context_type` is set to `LocalContext` (or its aliases `localcontext`, `Local`, `local`):

**remote\_profile:**

type: dict, optional  
 argument path: run\_mdata/model\_devi/  
 machine[LocalContext]/  
 remote\_profile

The information used to maintain the connection with remote machine. This field is empty for this context.

When `context_type` is set to `LazyLocalContext` (or its aliases `lazylocalcontext`, `LazyLocal`, `lazylocal`):

**remote\_profile:**

type: dict, optional  
 argument path: run\_mdata/model\_devi/  
 machine[LazyLocalContext]/  
 remote\_profile

The information used to maintain the connection with remote machine. This field is empty for this context.

When `context_type` is set to `SSHContext` (or its aliases `sshcontext`, `SSH`, `ssh`):

**remote\_profile:**

type: dict  
 argument path: run\_mdata/model\_devi/  
 machine[SSHContext]/remote\_profile

The information used to maintain the connection with remote machine.

**hostname:**

type: str  
 argument path:  
 run\_mdata/model\_devi/  
 machine[SSHContext]/  
 remote\_profile/hostname  
 hostname or ip of ssh connection.

**username:**

type: str  
 argument path:  
 run\_mdata/model\_devi/  
 machine[SSHContext]/  
 remote\_profile/username  
 username of target linux system

**password:**

type: str, optional

argument path:  
run\_mdata/model\_devi/  
machine[SSHContext]/  
remote\_profile/password

(deprecated) password of linux system. Please use [SSH keys](#) instead to improve security.

**port:**

type: int, optional, default: 22

argument path:  
run\_mdata/model\_devi/  
machine[SSHContext]/  
remote\_profile/port

ssh connection port.

**key\_filename:**

type: str | NoneType, optional,  
default: None

argument path:  
run\_mdata/model\_devi/  
machine[SSHContext]/  
remote\_profile/key\_filename

key filename used by ssh connection.  
If left None, find key in ~/.ssh or use  
password for login

**passphrase:**

type: str | NoneType, optional,  
default: None

argument path:  
run\_mdata/model\_devi/  
machine[SSHContext]/  
remote\_profile/passphrase

passphrase of key used by ssh connection

**timeout:**

type: int, optional, default: 10

argument path:  
run\_mdata/model\_devi/  
machine[SSHContext]/  
remote\_profile/timeout

timeout of ssh connection

**totp\_secret:**

type: str | NoneType, optional,  
default: None

argument path:  
run\_mdata/model\_devi/  
machine[SSHContext]/  
remote\_profile/totp\_secret

Time-based one time password secret. It should be a base32-encoded string extracted from the 2D code.

**tar\_compress:**

type: bool, optional, default: True

argument path:

run\_mdata/model\_devi/

machine[SSHContext]/

remote\_profile/tar\_compress

The archive will be compressed in upload and download if it is True. If not, compression will be skipped.

**look\_for\_keys:**

type: bool, optional, default: True

argument path:

run\_mdata/model\_devi/

machine[SSHContext]/

remote\_profile/look\_for\_keys

enable searching for discoverable private key files in ~/.ssh/

When `context_type` is set to `HDFSContext` (or its aliases `hdfscontext`, `HDFS`, `hdfs`):

**remote\_profile:**

type: dict, optional

argument path: run\_mdata/model\_devi/

machine[HDFSContext]/remote\_profile

The information used to maintain the connection with remote machine. This field is empty for this context.

When `context_type` is set to `OpenAPIContext` (or its aliases `openapicontext`, `OpenAPI`, `openapi`):

**remote\_profile:**

type: dict, optional

argument path: run\_mdata/model\_devi/

machine[OpenAPIContext]/

remote\_profile

The information used to maintain the connection with remote machine. This field is empty for this context.

**resources:**

type: dict

argument path: run\_mdata/model\_devi/resources

**number\_node:**

type: int, optional, default: 1

argument path: `run_mdata/model_devi/  
resources/number_node`

The number of node need for each *job*

**cpu\_per\_node:**

type: int, optional, default: 1

argument path: `run_mdata/model_devi/  
resources/cpu_per_node`

cpu numbers of each node assigned to each job.

**gpu\_per\_node:**

type: int, optional, default: 0

argument path: `run_mdata/model_devi/  
resources/gpu_per_node`

gpu numbers of each node assigned to each job.

**queue\_name:**

type: str, optional, default: (empty string)

argument path: `run_mdata/model_devi/  
resources/queue_name`

The queue name of batch job scheduler system.

**group\_size:**

type: int

argument path: `run_mdata/model_devi/  
resources/group_size`

The number of *tasks* in a *job*. 0 means infinity.

**custom\_flags:**

type: typing.List[str], optional

argument path: `run_mdata/model_devi/  
resources/custom_flags`

The extra lines pass to job submitting script  
header

**strategy:**

type: dict, optional

argument path: `run_mdata/model_devi/  
resources/strategy`

strategies we use to generation job submitting  
scripts.

**if\_cuda\_multi\_devices:**

type: bool, optional, default: False

argument path:  
`run_mdata/model_devi/  
resources/strategy/  
if_cuda_multi_devices`

If there are multiple nvidia GPUS on the node, and we want to assign the tasks to different GPUS. If true, dpdispatcher will manually export environment variable CUDA\_VISIBLE\_DEVICES to different task. Usually, this option will be used with Task.task\_need\_resources variable simultaneously.

**ratio\_unfinished:**

type: float, optional, default: 0.0  
 argument path: run\_mdata/  
 model\_devi/resources/  
 strategy/ratio\_unfinished

The ratio of *tasks* that can be unfinished.

**customized\_script\_header\_template\_file:**

type: str, optional  
 argument path:  
 run\_mdata/model\_devi/  
 resources/strategy/  
 customized\_script\_header\_template\_file

The customized template file to generate job submitting script header, which overrides the default file.

**para\_deg:**

type: int, optional, default: 1  
 argument path: run\_mdata/model\_devi/  
 resources/para\_deg

Decide how many tasks will be run in parallel.

**source\_list:**

type: typing.List[str], optional, default: []  
 argument path: run\_mdata/model\_devi/  
 resources/source\_list

The env file to be sourced before the command execution.

**module\_purge:**

type: bool, optional, default: False  
 argument path: run\_mdata/model\_devi/  
 resources/module\_purge

Remove all modules on HPC system before module load (module\_list)

**module\_unload\_list:**

type: `typing.List[str]`, optional, default: `[]`

argument path: `run_mdata/model_devi/resources/module_unload_list`

The modules to be unloaded on HPC system before submitting jobs

**module\_list:**

type: `typing.List[str]`, optional, default: `[]`

argument path: `run_mdata/model_devi/resources/module_list`

The modules to be loaded on HPC system before submitting jobs

**envs:**

type: `dict`, optional, default: `{}`

argument path: `run_mdata/model_devi/resources/envs`

The environment variables to be exported on before submitting jobs

**prepend\_script:**

type: `typing.List[str]`, optional, default: `[]`

argument path: `run_mdata/model_devi/resources/prepend_script`

Optional script run before jobs submitted.

**append\_script:**

type: `typing.List[str]`, optional, default: `[]`

argument path: `run_mdata/model_devi/resources/append_script`

Optional script run after jobs submitted.

**wait\_time:**

type: `float | int`, optional, default: `0`

argument path: `run_mdata/model_devi/resources/wait_time`

The waiting time in second after a single *task* submitted

Depending on the value of *batch\_type*, different sub args are accepted.



**batch\_type:**

type: str (flag key)

argument path: run\_mdata/model\_devi/  
resources/batch\_type

possible choices: *Bohrium*, *Fugaku*, *Slurm*,  
*SlurmJobArray*, *Shell*, *LSF*, *PBS*, *SGE*,  
*Torque*, *OpenAPI*, *DistributedShell*

The batch job system type loaded from machine/batch\_type.

When *batch\_type* is set to *Bohrium* (or its aliases *bohrium*, *Lebesgue*, *lebesgue*, *DpCloudServer*, *dpcloudserver*):

**kwargs:**

type: dict, optional

argument path: run\_mdata/model\_devi/  
resources[Bohrium]/kwargs

This field is empty for this batch.

When *batch\_type* is set to *Fugaku* (or its alias *fugaku*):

**kwargs:**

type: dict, optional

argument path: run\_mdata/model\_devi/  
resources[Fugaku]/kwargs

This field is empty for this batch.

When *batch\_type* is set to *Slurm* (or its alias *slurm*):

**kwargs:**

type: dict, optional

argument path: run\_mdata/model\_devi/  
resources[Slurm]/kwargs

Extra arguments.

**custom\_gpu\_line:**

type: str | NoneType, optional,

default: None

argument path: run\_mdata/  
model\_devi/resources[Slurm]/  
kwargs/custom\_gpu\_line

Custom GPU configuration, starting  
with #SBATCH

When *batch\_type* is set to *SlurmJobArray* (or its alias *slurmjobarray*):

**kwargs:**

type: dict, optional

argument path: `run_mdata/model_devi/  
resources[SlurmJobArray]/kwargs`

Extra arguments.

**custom\_gpu\_line:**

type: `str | NoneType`, optional,  
default: `None`

argument path:  
`run_mdata/model_devi/  
resources[SlurmJobArray]/  
kwargs/custom_gpu_line`

Custom GPU configuration, starting  
with `#SBATCH`

**slurm\_job\_size:**

type: `int`, optional, default: `1`

argument path:  
`run_mdata/model_devi/  
resources[SlurmJobArray]/  
kwargs/slurm_job_size`

Number of tasks in a Slurm job

When `batch_type` is set to `Shell` (or its alias `shell`):

**kwargs:**

type: `dict`, optional

argument path: `run_mdata/model_devi/  
resources[Shell]/kwargs`

This field is empty for this batch.

When `batch_type` is set to `LSF` (or its alias `lsf`):

**kwargs:**

type: `dict`

argument path: `run_mdata/model_devi/  
resources[LSF]/kwargs`

Extra arguments.

**gpu\_usage:**

type: `bool`, optional, default: `False`

argument path: `run_mdata/  
model_devi/resources[LSF]/  
kwargs/gpu_usage`

Choosing if GPU is used in the calcu-  
lation step.

**gpu\_new\_syntax:**

type: `bool`, optional, default: `False`

argument path: `run_mdata/  
model_devi/resources[LSF]/  
kwargs/gpu_new_syntax`

For LFS  $\geq$  10.1.0.3, new option -gpu for #BSUB could be used. If False, and old syntax would be used.

**gpu\_exclusive:**

type: bool, optional, default: True  
 argument path: run\_mdata/  
 model\_devi/resources[LSF]/  
 kwargs/gpu\_exclusive

Only take effect when new syntax enabled. Control whether submit tasks in exclusive way for GPU.

**custom\_gpu\_line:**

type: str | NoneType, optional,  
 default: None  
 argument path: run\_mdata/  
 model\_devi/resources[LSF]/  
 kwargs/custom\_gpu\_line

Custom GPU configuration, starting with #BSUB

When `batch_type` is set to PBS (or its alias pbs):

**kwargs:**

type: dict, optional  
 argument path: run\_mdata/model\_devi/  
 resources[PBS]/kwargs

This field is empty for this batch.

When `batch_type` is set to SGE (or its alias sge):

**kwargs:**

type: dict, optional  
 argument path: run\_mdata/model\_devi/  
 resources[SGE]/kwargs

This field is empty for this batch.

When `batch_type` is set to Torque (or its alias torque):

**kwargs:**

type: dict, optional  
 argument path: run\_mdata/model\_devi/  
 resources[Torque]/kwargs

This field is empty for this batch.

When `batch_type` is set to OpenAPI (or its alias openapi):

**kwargs:**

type: dict, optional

argument path: `run_mdata/model_devi/  
resources[OpenAPI]/kwargs`

This field is empty for this batch.

When `batch_type` is set to `DistributedShell` (or its alias `distributedshell`):

**kwargs:**

type: dict, optional

argument path: `run_mdata/model_devi/  
resources[DistributedShell]/kwargs`

This field is empty for this batch.

**user\_forward\_files:**

type: list, optional

argument path:

`run_mdata/model_devi/user_forward_files`

Files to be forwarded to the remote machine.

**user\_backward\_files:**

type: list, optional

argument path:

`run_mdata/model_devi/user_backward_files`

Files to be backwarded from the remote machine.

**fp:**

type: dict

argument path: `run_mdata/fp`

Parameters of command, machine, and resources for fp

**command:**

type: str

argument path: `run_mdata/fp/command`

Command of a program.

**machine:**

type: dict

argument path: `run_mdata/fp/machine`

**batch\_type:**

type: str

argument path:

`run_mdata/fp/machine/batch_type`

The batch job system type. Option: Fugaku, Slurm, DistributedShell, SGE, PBS, Torque, SlurmJobArray, Shell, Bohrium, LSF, OpenAPI

**local\_root:**

type: str | NoneType

argument path:

run\_mdata/fp/machine/local\_root

The dir where the tasks and relating files locate.  
Typically the project dir.

**remote\_root:**

type: str | NoneType, optional

argument path:

run\_mdata/fp/machine/remote\_root

The dir where the tasks are executed on the remote machine. Only needed when context is not lazy-local.

**clean\_asynchronously:**

type: bool, optional, default: False

argument path: run\_mdata/fp/machine/  
clean\_asynchronously

Clean the remote directory asynchronously after the job finishes.

Depending on the value of *context\_type*, different sub args are accepted.

**context\_type:**

type: str (flag key)

argument path:

run\_mdata/fp/machine/context\_type

possible choices: *BohriumContext*,  
*LocalContext*, *LazyLocalContext*,  
*SSHContext*, *HDFSContext*,  
*OpenAPIContext*

The connection used to remote machine. Option: OpenAPIContext, SSHContext, BohriumContext, HDFSContext, LazyLocalContext, LocalContext

When *context\_type* is set to *BohriumContext* (or its aliases *bohriumcontext*, *Bohrium*, *bohrium*, *DpCloudServerContext*, *dpcloudservercontext*, *DpCloudServer*, *dpcloudserver*, *LebesgueContext*, *lebesguecontext*, *Lebesgue*, *lebesgue*):

**remote\_profile:**

type: dict

argument path: run\_mdata/fp/  
machine[BohriumContext]/  
remote\_profile

The information used to maintain the connection with remote machine.

**email:**

type: str, optional  
argument path: run\_mdata/fp/  
machine[BohriumContext]/  
remote\_profile/email  
Email

**password:**

type: str, optional  
argument path: run\_mdata/fp/  
machine[BohriumContext]/  
remote\_profile/password  
Password

**program\_id:**

type: int, alias: *project\_id*  
argument path: run\_mdata/fp/  
machine[BohriumContext]/  
remote\_profile/program\_id  
Program ID

**retry\_count:**

type: NoneType | int, optional,  
default: 2  
argument path: run\_mdata/fp/  
machine[BohriumContext]/  
remote\_profile/retry\_count  
The retry count when a job is terminated

**ignore\_exit\_code:**

type: bool, optional, default: True  
argument path: run\_mdata/fp/  
machine[BohriumContext]/  
remote\_profile/  
ignore\_exit\_code

**The job state will be marked as finished if the exit code is non-zero when set to True. Otherwise,**  
the job state will be designated as terminated.

**keep\_backup:**

type: bool, optional  
argument path: run\_mdata/fp/  
machine[BohriumContext]/  
remote\_profile/keep\_backup  
keep download and upload zip

**input\_data:**

type: dict  
 argument path: run\_mdata/fp/  
 machine[BohriumContext]/  
 remote\_profile/input\_data

Configuration of job

When `context_type` is set to `LocalContext` (or its aliases `localcontext`, `Local`, `local`):

**remote\_profile:**

type: dict, optional  
 argument path:  
 run\_mdata/fp/machine[LocalContext]/  
 remote\_profile

The information used to maintain the connection with remote machine. This field is empty for this context.

When `context_type` is set to `LazyLocalContext` (or its aliases `lazylocalcontext`, `LazyLocal`, `lazylocal`):

**remote\_profile:**

type: dict, optional  
 argument path: run\_mdata/fp/  
 machine[LazyLocalContext]/  
 remote\_profile

The information used to maintain the connection with remote machine. This field is empty for this context.

When `context_type` is set to `SSHContext` (or its aliases `sshcontext`, `SSH`, `ssh`):

**remote\_profile:**

type: dict  
 argument path: run\_mdata/fp/  
 machine[SSHContext]/remote\_profile

The information used to maintain the connection with remote machine.

**hostname:**

type: str  
 argument path: run\_mdata/fp/  
 machine[SSHContext]/  
 remote\_profile/hostname

hostname or ip of ssh connection.

**username:**

type: str

argument path: `run_mdata/fp/machine[SSHContext]/remote_profile/username`

username of target linux system

**password:**

type: `str`, optional

argument path: `run_mdata/fp/machine[SSHContext]/remote_profile/password`

(deprecated) password of linux system. Please use [SSH keys](#) instead to improve security.

**port:**

type: `int`, optional, default: 22

argument path: `run_mdata/fp/machine[SSHContext]/remote_profile/port`

ssh connection port.

**key\_filename:**

type: `str | NoneType`, optional,  
default: `None`

argument path: `run_mdata/fp/machine[SSHContext]/remote_profile/key_filename`

key filename used by ssh connection.  
If left `None`, find key in `~/.ssh` or use  
password for login

**passphrase:**

type: `str | NoneType`, optional,  
default: `None`

argument path: `run_mdata/fp/machine[SSHContext]/remote_profile/passphrase`

passphrase of key used by ssh connection

**timeout:**

type: `int`, optional, default: 10

argument path: `run_mdata/fp/machine[SSHContext]/remote_profile/timeout`

timeout of ssh connection

**totp\_secret:**

type: `str | NoneType`, optional,  
default: `None`



argument path: `run_mdata/fp/machine[SSHContext]/remote_profile/totp_secret`

Time-based one time password secret. It should be a base32-encoded string extracted from the 2D code.

**tar\_compress:**

type: bool, optional, default: True  
 argument path: `run_mdata/fp/machine[SSHContext]/remote_profile/tar_compress`

The archive will be compressed in upload and download if it is True. If not, compression will be skipped.

**look\_for\_keys:**

type: bool, optional, default: True  
 argument path: `run_mdata/fp/machine[SSHContext]/remote_profile/look_for_keys`

enable searching for discoverable private key files in `~/.ssh/`

When `context_type` is set to `HDFSContext` (or its aliases `hdfscontext`, `HDFS`, `hdfs`):

**remote\_profile:**

type: dict, optional  
 argument path: `run_mdata/fp/machine[HDFSContext]/remote_profile`

The information used to maintain the connection with remote machine. This field is empty for this context.

When `context_type` is set to `OpenAPIContext` (or its aliases `openapicontext`, `OpenAPI`, `openapi`):

**remote\_profile:**

type: dict, optional  
 argument path: `run_mdata/fp/machine[OpenAPIContext]/remote_profile`

The information used to maintain the connection with remote machine. This field is empty for this context.

**resources:**

type: dict  
 argument path: `run_mdata/fp/resources`

**number\_node:**

type: int, optional, default: 1

argument path:

run\_mdata/fp/resources/number\_node

The number of node need for each *job*

**cpu\_per\_node:**

type: int, optional, default: 1

argument path:

run\_mdata/fp/resources/cpu\_per\_node

cpu numbers of each node assigned to each job.

**gpu\_per\_node:**

type: int, optional, default: 0

argument path:

run\_mdata/fp/resources/gpu\_per\_node

gpu numbers of each node assigned to each job.

**queue\_name:**

type: str, optional, default: (empty string)

argument path:

run\_mdata/fp/resources/queue\_name

The queue name of batch job scheduler system.

**group\_size:**

type: int

argument path:

run\_mdata/fp/resources/group\_size

The number of *tasks* in a *job*. 0 means infinity.

**custom\_flags:**

type: typing.List[str], optional

argument path:

run\_mdata/fp/resources/custom\_flags

The extra lines pass to job submitting script header

**strategy:**

type: dict, optional

argument path:

run\_mdata/fp/resources/strategy

strategies we use to generation job submitting scripts.

**if\_cuda\_multi\_devices:**

type: bool, optional, default: False

argument path: `run_mdata/fp/  
resources/strategy/  
if_cuda_multi_devices`

If there are multiple nvidia GPUS on the node, and we want to assign the tasks to different GPUS. If true, `dpdispatcher` will manually export environment variable `CUDA_VISIBLE_DEVICES` to different task. Usually, this option will be used with `Task.task_need_resources` variable simultaneously.

#### **ratio\_unfinished:**

type: float, optional, default: 0.0

argument path:

`run_mdata/fp/resources/  
strategy/ratio_unfinished`

The ratio of *tasks* that can be unfinished.

#### **customized\_script\_header\_template\_file:**

type: str, optional

argument path: `run_mdata/fp/  
resources/strategy/  
customized_script_header_template_file`

The customized template file to generate job submitting script header, which overrides the default file.

#### **para\_deg:**

type: int, optional, default: 1

argument path:

`run_mdata/fp/resources/para_deg`

Decide how many tasks will be run in parallel.

#### **source\_list:**

type: `typing.List[str]`, optional, default: []

argument path:

`run_mdata/fp/resources/source_list`

The env file to be sourced before the command execution.

#### **module\_purge:**

type: bool, optional, default: False

argument path:

`run_mdata/fp/resources/module_purge`

Remove all modules on HPC system before module load (`module_list`)

**module\_unload\_list:**

type: `typing.List[str]`, optional, default: `[]`

argument path: `run_mdata/fp/resources/module_unload_list`

The modules to be unloaded on HPC system before submitting jobs

**module\_list:**

type: `typing.List[str]`, optional, default: `[]`

argument path: `run_mdata/fp/resources/module_list`

The modules to be loaded on HPC system before submitting jobs

**envs:**

type: `dict`, optional, default: `{}`

argument path: `run_mdata/fp/resources/envs`

The environment variables to be exported on before submitting jobs

**prepend\_script:**

type: `typing.List[str]`, optional, default: `[]`

argument path: `run_mdata/fp/resources/prepend_script`

Optional script run before jobs submitted.

**append\_script:**

type: `typing.List[str]`, optional, default: `[]`

argument path: `run_mdata/fp/resources/append_script`

Optional script run after jobs submitted.

**wait\_time:**

type: `float | int`, optional, default: `0`

argument path: `run_mdata/fp/resources/wait_time`

The waiting time in second after a single *task* submitted

Depending on the value of *batch\_type*, different sub args are accepted.

**batch\_type:**

type: str (flag key)

argument path:

run\_mdata/fp/resources/batch\_type

possible choices: *Bohrium*, *Fugaku*, *Slurm*,  
*SlurmJobArray*, *Shell*, *LSF*, *PBS*, *SGE*,  
*Torque*, *OpenAPI*, *DistributedShell*

The batch job system type loaded from machine/batch\_type.

When *batch\_type* is set to *Bohrium* (or its aliases *bohrium*, *Lebesgue*, *lebesgue*, *DpCloudServer*, *dpcloudserver*):

**kwargs:**

type: dict, optional

argument path: run\_mdata/fp/  
resources[Bohrium]/kwargs

This field is empty for this batch.

When *batch\_type* is set to *Fugaku* (or its alias *fugaku*):

**kwargs:**

type: dict, optional

argument path: run\_mdata/fp/  
resources[Fugaku]/kwargs

This field is empty for this batch.

When *batch\_type* is set to *Slurm* (or its alias *slurm*):

**kwargs:**

type: dict, optional

argument path: run\_mdata/fp/  
resources[Slurm]/kwargs

Extra arguments.

**custom\_gpu\_line:**

type: str | NoneType, optional,  
default: None

argument path: run\_mdata/fp/  
resources[Slurm]/kwargs/  
custom\_gpu\_line

Custom GPU configuration, starting  
with #SBATCH

When *batch\_type* is set to *SlurmJobArray* (or its alias *slurmjobarray*):

**kwargs:**

type: dict, optional

argument path: `run_mdata/fp/resources[SlurmJobArray]/kwargs`

Extra arguments.

**custom\_gpu\_line:**

type: `str | NoneType`, optional,  
default: `None`

argument path: `run_mdata/fp/resources[SlurmJobArray]/kwargs/custom_gpu_line`

Custom GPU configuration, starting with `#SBATCH`

**slurm\_job\_size:**

type: `int`, optional, default: `1`  
argument path: `run_mdata/fp/resources[SlurmJobArray]/kwargs/slurm_job_size`

Number of tasks in a Slurm job

When `batch_type` is set to `Shell` (or its alias `shell`):

**kwargs:**

type: `dict`, optional  
argument path: `run_mdata/fp/resources[Shell]/kwargs`

This field is empty for this batch.

When `batch_type` is set to `LSF` (or its alias `lsf`):

**kwargs:**

type: `dict`  
argument path: `run_mdata/fp/resources[LSF]/kwargs`

Extra arguments.

**gpu\_usage:**

type: `bool`, optional, default: `False`  
argument path: `run_mdata/fp/resources[LSF]/kwargs/gpu_usage`

Choosing if GPU is used in the calculation step.

**gpu\_new\_syntax:**

type: `bool`, optional, default: `False`  
argument path: `run_mdata/fp/resources[LSF]/kwargs/gpu_new_syntax`

For LFS  $\geq$  10.1.0.3, new option `-gpu` for `#BSUB` could be used. If False, and old syntax would be used.

**gpu\_exclusive:**

type: bool, optional, default: True  
 argument path:  
`run_mdata/fp/resources[LSF]/kwargs/gpu_exclusive`

Only take effect when new syntax enabled. Control whether submit tasks in exclusive way for GPU.

**custom\_gpu\_line:**

type: str | NoneType, optional,  
 default: None  
 argument path:  
`run_mdata/fp/resources[LSF]/kwargs/custom_gpu_line`

Custom GPU configuration, starting with `#BSUB`

When `batch_type` is set to PBS (or its alias `pbs`):

**kwargs:**

type: dict, optional  
 argument path:  
`run_mdata/fp/resources[PBS]/kwargs`

This field is empty for this batch.

When `batch_type` is set to SGE (or its alias `sge`):

**kwargs:**

type: dict, optional  
 argument path:  
`run_mdata/fp/resources[SGE]/kwargs`

This field is empty for this batch.

When `batch_type` is set to Torque (or its alias `torque`):

**kwargs:**

type: dict, optional  
 argument path: `run_mdata/fp/resources[Torque]/kwargs`

This field is empty for this batch.

When `batch_type` is set to OpenAPI (or its alias `openapi`):

**kwargs:**

type: dict, optional

argument path: `run_mdata/fp/`  
`resources[OpenAPI]/kwargs`

This field is empty for this batch.

When `batch_type` is set to `DistributedShell` (or its alias `distributedshell`):

**kwargs:**

type: dict, optional  
argument path: `run_mdata/fp/`  
`resources[DistributedShell]/kwargs`

This field is empty for this batch.

**user\_forward\_files:**

type: list, optional  
argument path: `run_mdata/fp/user_forward_files`

Files to be forwarded to the remote machine.

**user\_backward\_files:**

type: list, optional  
argument path: `run_mdata/fp/user_backward_files`

Files to be backwarded from the remote machine.



## 5.1 Init\_bulk

You may prepare initial data for bulk systems with VASP by:

```
dpngen init_bulk PARAM [MACHINE]
```

The MACHINE configure file is optional. If this parameter exists, then the optimization tasks or MD tasks will be submitted automatically according to MACHINE.json.

Basically init\_bulk can be divided into four parts , denoted as stages in PARAM:

1. Relax in folder 00.place\_ele
2. Perturb and scale in folder 01.scale\_pert
3. Run a short AIMD in folder 02.md
4. Collect data in folder 02.md.

All stages must be **in order**. One doesn't need to run all stages. For example, you may run stage 1 and 2, generating supercells as starting point of exploration in dpngen run.

If MACHINE is None, there should be only one stage in stages. Corresponding tasks will be generated, but user's intervention should be involved in, to manually run the scripts.

Following is an example for PARAM, which generates data from a typical structure hcp.

```
{
  "stages" : [1,2,3,4],
  "cell_type": "hcp",
  "latt": 4.479,
  "super_cell": [2, 2, 2],
  "elements": ["Mg"],
  "potcars": ["...../POTCAR"],
  "relax_incar": "...../INCAR_metal_rlx",
  "md_incar" : "...../INCAR_metal_md",
  "scale": [1.00],
  "skip_relax": false,
  "pert_numb": 2,
  "md_nstep" : 5,
  "pert_box": 0.03,
  "pert_atom": 0.01,
  "coll_ndata": 5000,
  "type_map" : [ "Mg", "Al"],
}
```

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```

    "_comment":      "that's all"
}

```

If you want to specify a structure as starting point for `init_bulk`, you may set in `PARAM` as follows.

```

"from_poscar":      true,
"from_poscar_path": "...../C_mp-47_conventional.POSCAR",

```

`init_bulk` supports both VASP and ABACUS for first-principle calculation. You can choose the software by specifying the key `init_fp_style`. If `init_fp_style` is not specified, the default software will be VASP.

When using ABACUS for `init_fp_style`, the keys of the paths of INPUT files for relaxation and MD simulations are the same as INCAR for VASP, which are `relax_incar` and `md_incar` respectively. Use `relax_kpt` and `md_kpt` for the relative path for KPT files of relaxation and MD simulations. They two can be omitted if `kspacing` (in unit of 1/Bohr) or `gamma_only` has been set in corresponding INPUT files. If `from_poscar` is set to `false`, you have to specify `atom_masses` in the same order as elements.

## 5.2 dpngen init\_bulk parameters

**Note:** One can load, modify, and export the input file by using our effective web-based tool [DP-GUI](#) online or hosted using the *command line interface* `dpngen gui`. All parameters below can be set in DP-GUI. By clicking “SAVE JSON”, one can download the input file.

### `init_bulk_jdata:`

type: dict

argument path: `init_bulk_jdata`

Generate initial data for bulk systems.

#### `stages:`

type: list[int]

argument path: `init_bulk_jdata/stages`

Stages for *init\_bulk*.

#### `elements:`

type: list[str]

argument path: `init_bulk_jdata/elements`

Atom types.

#### `potcars:`

type: list[str], optional

argument path: `init_bulk_jdata/potcars`

Path of POTCAR.

#### `cell_type:`

type: str, optional

argument path: `init_bulk_jdata/cell_type`

Specifying which typical structure to be generated. **Options** include `fcc`, `hcp`, `bcc`, `sc`, `diamond`.

**super\_cell:**

type: list[int]  
 argument path: init\_bulk\_jdata/super\_cell  
 Size of supercell.

**from\_poscar:**

type: bool, optional, default: False  
 argument path: init\_bulk\_jdata/from\_poscar  
 Deciding whether to use a given poscar as the beginning of relaxation. If it's true, keys (*cell\_type*, *latt*) will be aborted. Otherwise, these two keys are **necessary**.

**from\_poscar\_path:**

type: str, optional  
 argument path: init\_bulk\_jdata/from\_poscar\_path  
 Path of POSCAR for VASP or STRU for ABACUS. **Necessary** if *from\_poscar* is true.

**relax\_incar:**

type: str, optional  
 argument path: init\_bulk\_jdata/relax\_incar  
 Path of INCAR for VASP or INPUT for ABACUS for relaxation in VASP. **Necessary** if *stages* include 1.

**md\_incar:**

type: str, optional  
 argument path: init\_bulk\_jdata/md\_incar  
 Path of INCAR for VASP or INPUT for ABACUS for MD in VASP. **Necessary** if *stages* include 3.

**scale:**

type: list[float]  
 argument path: init\_bulk\_jdata/scale  
 Scales for isotropic transforming cells.

**skip\_relax:**

type: bool  
 argument path: init\_bulk\_jdata/skip\_relax  
 If it's true, you may directly run stage 2 (perturb and scale) using an unrelaxed POSCAR.

**pert\_numb:**

type: int  
 argument path: init\_bulk\_jdata/pert\_numb  
 Number of perturbations for each scaled (key *scale*) POSCAR.

**pert\_box:**

type: float  
 argument path: init\_bulk\_jdata/pert\_box

Anisotropic Perturbation for cells (independent changes of lengths of three box vectors as well as angle among) in decimal formats. 9 elements of the 3x3 perturbation matrix will be randomly sampled from a uniform distribution (default) in the range  $[-\text{pert\_box}, \text{pert\_box}]$ . Such a perturbation matrix adds the identity matrix gives the actual transformation matrix for this perturbation operation.

**pert\_atom:**

type: float

argument path: `init_bulk_jdata/pert_atom`

Perturbation of atom coordinates (Angstrom). Random perturbations are performed on three coordinates of each atom by adding values randomly sampled from a uniform distribution in the range  $[-\text{pert\_atom}, \text{pert\_atom}]$ .

**md\_nstep:**

type: int

argument path: `init_bulk_jdata/md_nstep`

Steps of AIMD in stage 3. If it's not equal to settings via *NSW* in *md\_incar*, DP-GEN will follow *NSW*.

**coll\_ndata:**

type: int

argument path: `init_bulk_jdata/coll_ndata`

Maximal number of collected data.

**type\_map:**

type: `list[str]`, optional

argument path: `init_bulk_jdata/type_map`

The indices of elements in deepmd formats will be set in this order.

Depending on the value of *init\_fp\_style*, different sub args are accepted.

**init\_fp\_style:**

type: `str` (flag key), default: VASP

argument path: `init_bulk_jdata/init_fp_style`

possible choices: [VASP](#), [ABACUS](#)

First-principle software. If this key is absent.

When *init\_fp\_style* is set to VASP:

No more parameters is needed to be added.

When *init\_fp\_style* is set to ABACUS:

ABACUS

**relax\_kpt:**

type: `str`, optional

argument path: `init_bulk_jdata[ABACUS]/relax_kpt`

Path of *KPT* file for relaxation in stage 1. Only useful if *init\_fp\_style* is "ABACUS".

**md\_kpt:**

type: `str`, optional

argument path: `init_bulk_jdata[ABACUS]/md_kpt`

Path of *KPT* file for MD simulations in stage 3. Only useful if *init\_fp\_style* is “ABACUS”.

**atom\_masses:**

type: `list[float]`, optional

argument path: `init_bulk_jdata[ABACUS]/atom_masses`

List of atomic masses of elements. The order should be the same as *Elements*.

Only useful if *init\_fp\_style* is “ABACUS”.

## 5.3 dpngen init\_bulk machine parameters

---

**Note:** One can load, modify, and export the input file by using our effective web-based tool [DP-GUI](#) online or hosted using the *command line interface* `dpngen gui`. All parameters below can be set in DP-GUI. By clicking “SAVE JSON”, one can download the input file.

---

**init\_bulk\_mdata:**

type: `dict`

argument path: `init_bulk_mdata`

machine.json file

**api\_version:**

type: `str`, optional, default: 1.0

argument path: `init_bulk_mdata/api_version`

Please set to 1.0

**deepmd\_version:**

type: `str`, optional, default: 2

argument path: `init_bulk_mdata/deepmd_version`

DeePMD-kit version, e.g. 2.1.3

**fp:**

type: `dict`

argument path: `init_bulk_mdata/fp`

Parameters of command, machine, and resources for fp

**command:**

type: `str`

argument path: `init_bulk_mdata/fp/command`

Command of a program.

**machine:**

type: `dict`

argument path: `init_bulk_mdata/fp/machine`

**batch\_type:**

type: str  
argument path: init\_bulk\_mdata/fp/  
machine/batch\_type

The batch job system type. Option: Fugaku, Slurm, DistributedShell, SGE, PBS, Torque, SlurmJobArray, Shell, Bohrium, LSF, OpenAPI

**local\_root:**

type: str | NoneType  
argument path: init\_bulk\_mdata/fp/  
machine/local\_root

The dir where the tasks and relating files locate. Typically the project dir.

**remote\_root:**

type: str | NoneType, optional  
argument path: init\_bulk\_mdata/fp/  
machine/remote\_root

The dir where the tasks are executed on the remote machine. Only needed when context is not lazy-local.

**clean\_asynchronously:**

type: bool, optional, default: False  
argument path: init\_bulk\_mdata/fp/  
machine/clean\_asynchronously

Clean the remote directory asynchronously after the job finishes.

Depending on the value of *context\_type*, different sub args are accepted.

**context\_type:**

type: str (flag key)  
argument path: init\_bulk\_mdata/fp/  
machine/context\_type  
possible choices: *BohriumContext*,  
*LocalContext*, *LazyLocalContext*,  
*SSHContext*, *HDFSContext*,  
*OpenAPIContext*

The connection used to remote machine. Option: OpenAPIContext, SSHContext, BohriumContext, HDFSContext, LazyLocalContext, LocalContext

When *context\_type* is set to *BohriumContext* (or its aliases *bohriumcontext*, *Bohrium*, *bohrium*, *DpCloudServerContext*, *dpcloudservercontext*,

DpCloudServer, dpcloudserver, LebesgueContext, lebesguecontext, Lebesgue, lebesgue):

**remote\_profile:**

type: dict  
 argument path: init\_bulk\_mdata/fp/  
 machine[BohriumContext]/  
 remote\_profile

The information used to maintain the connection with remote machine.

**email:**

type: str, optional  
 argument path: init\_bulk\_mdata/  
 fp/machine[BohriumContext]/  
 remote\_profile/email

Email

**password:**

type: str, optional  
 argument path: init\_bulk\_mdata/  
 fp/machine[BohriumContext]/  
 remote\_profile/password

Password

**program\_id:**

type: int, alias: *project\_id*  
 argument path: init\_bulk\_mdata/  
 fp/machine[BohriumContext]/  
 remote\_profile/program\_id

Program ID

**retry\_count:**

type: NoneType | int, optional,  
 default: 2  
 argument path: init\_bulk\_mdata/  
 fp/machine[BohriumContext]/  
 remote\_profile/retry\_count

The retry count when a job is terminated

**ignore\_exit\_code:**

type: bool, optional, default: True  
 argument path: init\_bulk\_mdata/  
 fp/machine[BohriumContext]/  
 remote\_profile/  
 ignore\_exit\_code

**The job state will be marked as finished if the exit code is non-zero when set to True. Otherwise, the job state will be designated as terminated.**

**keep\_backup:**

type: bool, optional  
argument path: init\_bulk\_mdata/  
fp/machine[BohriumContext]/  
remote\_profile/keep\_backup  
keep download and upload zip

**input\_data:**

type: dict  
argument path: init\_bulk\_mdata/  
fp/machine[BohriumContext]/  
remote\_profile/input\_data  
Configuration of job

When `context_type` is set to `LocalContext` (or its aliases `localcontext`, `Local`, `local`):

**remote\_profile:**

type: dict, optional  
argument path: init\_bulk\_mdata/fp/  
machine[LocalContext]/  
remote\_profile  
The information used to maintain the connection with remote machine. This field is empty for this context.

When `context_type` is set to `LazyLocalContext` (or its aliases `lazylocalcontext`, `LazyLocal`, `lazylocal`):

**remote\_profile:**

type: dict, optional  
argument path: init\_bulk\_mdata/fp/  
machine[LazyLocalContext]/  
remote\_profile  
The information used to maintain the connection with remote machine. This field is empty for this context.

When `context_type` is set to `SSHContext` (or its aliases `sshcontext`, `SSH`, `ssh`):

**remote\_profile:**

type: dict  
argument path: init\_bulk\_mdata/fp/  
machine[SSHContext]/remote\_profile  
The information used to maintain the connection with remote machine.

**hostname:**

type: str



argument path: `init_bulk_mdata/  
fp/machine[SSHContext]/  
remote_profile/hostname`

hostname or ip of ssh connection.

**username:**

type: `str`

argument path: `init_bulk_mdata/  
fp/machine[SSHContext]/  
remote_profile/username`

username of target linux system

**password:**

type: `str`, optional

argument path: `init_bulk_mdata/  
fp/machine[SSHContext]/  
remote_profile/password`

(deprecated) password of linux system. Please use [SSH keys](#) instead to improve security.

**port:**

type: `int`, optional, default: 22

argument path: `init_bulk_mdata/  
fp/machine[SSHContext]/  
remote_profile/port`

ssh connection port.

**key\_filename:**

type: `str | NoneType`, optional,  
default: `None`

argument path: `init_bulk_mdata/  
fp/machine[SSHContext]/  
remote_profile/key_filename`

key filename used by ssh connection.  
If left `None`, find key in `~/.ssh` or use  
password for login

**passphrase:**

type: `str | NoneType`, optional,  
default: `None`

argument path: `init_bulk_mdata/  
fp/machine[SSHContext]/  
remote_profile/passphrase`

passphrase of key used by ssh connection

**timeout:**

type: `int`, optional, default: 10

argument path: `init_bulk_mdata/  
fp/machine[SSHContext]/  
remote_profile/timeout`

timeout of ssh connection

**totp\_secret:**

type: str | NoneType, optional,  
default: None

argument path: init\_bulk\_mdata/  
fp/machine[SSHContext]/  
remote\_profile/totp\_secret

Time-based one time password secret. It should be a base32-encoded string extracted from the 2D code.

**tar\_compress:**

type: bool, optional, default: True  
argument path: init\_bulk\_mdata/  
fp/machine[SSHContext]/  
remote\_profile/tar\_compress

The archive will be compressed in upload and download if it is True. If not, compression will be skipped.

**look\_for\_keys:**

type: bool, optional, default: True  
argument path: init\_bulk\_mdata/  
fp/machine[SSHContext]/  
remote\_profile/look\_for\_keys

enable searching for discoverable private key files in ~/.ssh/

When `context_type` is set to `HDFSContext` (or its aliases `hdfscontext`, `HDFS`, `hdfs`):

**remote\_profile:**

type: dict, optional  
argument path: init\_bulk\_mdata/fp/  
machine[HDFSContext]/remote\_profile

The information used to maintain the connection with remote machine. This field is empty for this context.

When `context_type` is set to `OpenAPIContext` (or its aliases `openapicontext`, `OpenAPI`, `openapi`):

**remote\_profile:**

type: dict, optional  
argument path: init\_bulk\_mdata/fp/  
machine[OpenAPIContext]/  
remote\_profile

The information used to maintain the connection with remote machine. This field is empty for this context.

**resources:**

type: dict

argument path: `init_bulk_mdata/fp/resources`

**number\_node:**

type: int, optional, default: 1

argument path: `init_bulk_mdata/fp/resources/number_node`

The number of node need for each *job*

**cpu\_per\_node:**

type: int, optional, default: 1

argument path: `init_bulk_mdata/fp/resources/cpu_per_node`

cpu numbers of each node assigned to each job.

**gpu\_per\_node:**

type: int, optional, default: 0

argument path: `init_bulk_mdata/fp/resources/gpu_per_node`

gpu numbers of each node assigned to each job.

**queue\_name:**

type: str, optional, default: (empty string)

argument path: `init_bulk_mdata/fp/resources/queue_name`

The queue name of batch job scheduler system.

**group\_size:**

type: int

argument path: `init_bulk_mdata/fp/resources/group_size`

The number of *tasks* in a *job*. 0 means infinity.

**custom\_flags:**

type: typing.List[str], optional

argument path: `init_bulk_mdata/fp/resources/custom_flags`

The extra lines pass to job submitting script header

**strategy:**

type: dict, optional

argument path: `init_bulk_mdata/fp/resources/strategy`

strategies we use to generation job submitting scripts.

**if\_cuda\_multi\_devices:**

type: bool, optional, default: False  
argument path: `init_bulk_mdata/  
fp/resources/strategy/  
if_cuda_multi_devices`

If there are multiple nvidia GPUS on the node, and we want to assign the tasks to different GPUS. If true, `dpdispatcher` will manually export environment variable `CUDA_VISIBLE_DEVICES` to different task. Usually, this option will be used with `Task.task_need_resources` variable simultaneously.

**ratio\_unfinished:**

type: float, optional, default: 0.0  
argument path: `init_bulk_mdata/  
fp/resources/strategy/  
ratio_unfinished`

The ratio of *tasks* that can be unfinished.

**customized\_script\_header\_template\_file:**

type: str, optional  
argument path: `init_bulk_mdata/  
fp/resources/strategy/  
customized_script_header_template_file`

The customized template file to generate job submitting script header, which overrides the default file.

**para\_deg:**

type: int, optional, default: 1  
argument path: `init_bulk_mdata/fp/  
resources/para_deg`

Decide how many tasks will be run in parallel.

**source\_list:**

type: `typing.List[str]`, optional, default: []  
argument path: `init_bulk_mdata/fp/  
resources/source_list`

The env file to be sourced before the command execution.

**module\_purge:**

type: bool, optional, default: False  
argument path: `init_bulk_mdata/fp/  
resources/module_purge`

Remove all modules on HPC system before module load (module\_list)

**module\_unload\_list:**

type: `typing.List[str]`, optional, default: `[]`

argument path: `init_bulk_mdata/fp/resources/module_unload_list`

The modules to be unloaded on HPC system before submitting jobs

**module\_list:**

type: `typing.List[str]`, optional, default: `[]`

argument path: `init_bulk_mdata/fp/resources/module_list`

The modules to be loaded on HPC system before submitting jobs

**envs:**

type: `dict`, optional, default: `{}`

argument path: `init_bulk_mdata/fp/resources/envs`

The environment variables to be exported on before submitting jobs

**prepend\_script:**

type: `typing.List[str]`, optional, default: `[]`

argument path: `init_bulk_mdata/fp/resources/prepend_script`

Optional script run before jobs submitted.

**append\_script:**

type: `typing.List[str]`, optional, default: `[]`

argument path: `init_bulk_mdata/fp/resources/append_script`

Optional script run after jobs submitted.

**wait\_time:**

type: `float | int`, optional, default: `0`

argument path: `init_bulk_mdata/fp/resources/wait_time`

The waiting time in second after a single *task* submitted

Depending on the value of *batch\_type*, different sub args are accepted.

**batch\_type:**

type: str (flag key)  
argument path: `init_bulk_mdata/fp/resources/batch_type`  
possible choices: *Bohrium*, *Fugaku*, *Slurm*,  
*SlurmJobArray*, *Shell*, *LSF*, *PBS*, *SGE*,  
*Torque*, *OpenAPI*, *DistributedShell*  
The batch job system type loaded from machine/`batch_type`.

When `batch_type` is set to *Bohrium* (or its aliases *bohrium*, *Lebesgue*, *lebesgue*, *DpCloudServer*, *dpcloudserver*):

**kwargs:**

type: dict, optional  
argument path: `init_bulk_mdata/fp/resources[Bohrium]/kwargs`  
This field is empty for this batch.

When `batch_type` is set to *Fugaku* (or its alias *fugaku*):

**kwargs:**

type: dict, optional  
argument path: `init_bulk_mdata/fp/resources[Fugaku]/kwargs`  
This field is empty for this batch.

When `batch_type` is set to *Slurm* (or its alias *slurm*):

**kwargs:**

type: dict, optional  
argument path: `init_bulk_mdata/fp/resources[Slurm]/kwargs`  
Extra arguments.

**custom\_gpu\_line:**

type: str | NoneType, optional,  
default: None  
argument path: `init_bulk_mdata/fp/resources[Slurm]/kwargs/custom_gpu_line`  
Custom GPU configuration, starting with `#SBATCH`

When `batch_type` is set to *SlurmJobArray* (or its alias *slurmjobarray*):

**kwargs:**

type: dict, optional

argument path: `init_bulk_mdata/fp/resources[SlurmJobArray]/kwargs`

Extra arguments.

**custom\_gpu\_line:**

type: `str | NoneType`, optional,  
default: `None`

argument path: `init_bulk_mdata/fp/resources[SlurmJobArray]/kwargs/custom_gpu_line`

Custom GPU configuration, starting with `#SBATCH`

**slurm\_job\_size:**

type: `int`, optional, default: `1`

argument path: `init_bulk_mdata/fp/resources[SlurmJobArray]/kwargs/slurm_job_size`

Number of tasks in a Slurm job

When `batch_type` is set to `Shell` (or its alias `shell`):

**kwargs:**

type: `dict`, optional

argument path: `init_bulk_mdata/fp/resources[Shell]/kwargs`

This field is empty for this batch.

When `batch_type` is set to `LSF` (or its alias `lsf`):

**kwargs:**

type: `dict`

argument path: `init_bulk_mdata/fp/resources[LSF]/kwargs`

Extra arguments.

**gpu\_usage:**

type: `bool`, optional, default: `False`

argument path: `init_bulk_mdata/fp/resources[LSF]/kwargs/gpu_usage`

Choosing if GPU is used in the calculation step.

**gpu\_new\_syntax:**

type: `bool`, optional, default: `False`

argument path: `init_bulk_mdata/fp/resources[LSF]/kwargs/gpu_new_syntax`

For LFS  $\geq$  10.1.0.3, new option `-gpu` for `#BSUB` could be used. If False, and old syntax would be used.

**gpu\_exclusive:**

type: bool, optional, default: True  
argument path: `init_bulk_mdata/  
fp/resources[LSF]/kwargs/  
gpu_exclusive`

Only take effect when new syntax enabled. Control whether submit tasks in exclusive way for GPU.

**custom\_gpu\_line:**

type: str | NoneType, optional,  
default: None  
argument path: `init_bulk_mdata/  
fp/resources[LSF]/kwargs/  
custom_gpu_line`

Custom GPU configuration, starting with `#BSUB`

When `batch_type` is set to PBS (or its alias `pbs`):

**kwargs:**

type: dict, optional  
argument path: `init_bulk_mdata/fp/  
resources[PBS]/kwargs`

This field is empty for this batch.

When `batch_type` is set to SGE (or its alias `sge`):

**kwargs:**

type: dict, optional  
argument path: `init_bulk_mdata/fp/  
resources[SGE]/kwargs`

This field is empty for this batch.

When `batch_type` is set to Torque (or its alias `torque`):

**kwargs:**

type: dict, optional  
argument path: `init_bulk_mdata/fp/  
resources[Torque]/kwargs`

This field is empty for this batch.

When `batch_type` is set to OpenAPI (or its alias `openapi`):

**kwargs:**

type: dict, optional



argument path: `init_bulk_mdata/fp/resources[OpenAPI]/kwargs`

This field is empty for this batch.

When `batch_type` is set to `DistributedShell` (or its alias `distributedshell`):

**kwargs:**

type: dict, optional

argument path: `init_bulk_mdata/fp/resources[DistributedShell]/kwargs`

This field is empty for this batch.

**user\_forward\_files:**

type: list, optional

argument path:

`init_bulk_mdata/fp/user_forward_files`

Files to be forwarded to the remote machine.

**user\_backward\_files:**

type: list, optional

argument path:

`init_bulk_mdata/fp/user_backward_files`

Files to be backwarded from the remote machine.

## 5.4 Init\_surf

You may prepare initial data for surface systems with VASP by:

```
dpngen init_surf PARAM [MACHINE]
```

The MACHINE configure file is optional. If this parameter exists, then the optimization tasks or MD tasks will be submitted automatically according to MACHINE.json. That is to say, if one only wants to prepare `surf-xxx/sys-xxx` folders for the second stage but wants to skip relaxation, `dpngen init_surf PARAM` should be used (without MACHINE). “stages” and “skip\_relax” in PARAM should be set as:

```
"stages": [1,2],
"skip_relax": true,
```

Basically `init_surf` can be divided into two parts, denoted as `stages` in PARAM:

1. Build specific surface in folder `00.place_ele`
2. Perturb and scale in folder `01.scale_pert`

All stages must be **in order**.

Generally, `init_surf` does not run AIMD but only generates a lot of configurations. Compared with `init_bulk`, which runs DFT calculations twice, `init_surf` does once. Usually, we do `init_bulk`, run many rounds of DP-GEN iterations, collect enough data for the bulk system, and do `init_surf` after that. At this point, the lattice constant has been determined, and the lattice constant required for the initial configuration of `init_surf` can be used directly.

These configurations made by `init_surf` are prepared for `01.model_devi`. Candidates will do DFT calculation in `02.fp`.

- Generate vacuum layers

According to [the source code of `pert\_scaled`](#), `init_surf` will generate a series of surface structures with specified separations between the sample layer and its periodic image. There are two ways to specify the interval in generating the vacuum layers: 1) to set the interval value and 2) to set the number of intervals.

You can use `layer_numb` (the number of layers of the slab) or `z_min` (the total thickness) to specify the thickness of the atoms below. Then `vacuum_*` parameters specify the vacuum layers above. `dpgen init_surf` will make a series of structures with the thickness of vacuum layers from `vacuum_min` to `vacuum_max`. The number of vacuum layers is controlled by the parameter `vacuum_resol`.

The layers will be generated even when the size of `vacuum_resol` is 1. When the size of `vacuum_resol` is 2 or it is empty, the whole interval range is divided into the nearby region with denser intervals (head region) and the far-away region with sparser intervals (tail region), which are divided by `mid_point`.

When the size of `vacuum_resol` is 2, two elements respectively decide the number of intervals in head region and tail region.

When `vacuum_resol` is empty, the number of intervals in the head region = `vacuum_num` \* `head_ratio`. `vacuum_num` and `head_ratio` are both keys in `param.json`.

- Attach files in the task path

One can use the machine parameter `forward_files` to upload other files besides POSCAR, INCAR, and POTCAR. For example, “`vdw_kernel.bindat`” for each task.

See [the document of task parameters](#).

Following is an example for `PARAM`, which generates data from a typical structure `fcc`.

```
{
  "stages": [
    1,
    2
  ],
  "cell_type": "fcc",
  "latt": 4.034,
  "super_cell": [
    2,
    2,
    2
  ],
  "layer_numb": 3,
  "vacuum_max": 9.0,
  "vacuum_resol": [
    0.5,
    1
  ],
  "mid_point": 4.0,
  "millers": [
    [
      1,
      0,
      0
    ]
  ],
}
```

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```

[
  1,
  1,
  0
],
[
  1,
  1,
  1
]
],
"elements": [
  "Al"
],
"potcars": [
  "...../POTCAR"
],
"relax_incar": "...../INCAR_metal_rlx_low",
"scale": [
  1.0
],
"skip_relax": true,
"pert_numb": 2,
"pert_box": 0.03,
"pert_atom": 0.01,
"_comment": "that's all"
}

```

Another example is `from_poscar` method. Here you need to specify the POSCAR file.

```

{
  "stages": [
    1,
    2
  ],
  "cell_type": "fcc",
  "from_poscar": true,
  "from_poscar_path": "POSCAR",
  "super_cell": [
    1,
    1,
    1
  ],
  "layer_numb": 3,
  "vacuum_max": 5.0,
  "vacuum_resol": [0.5, 2],
  "mid_point": 2.0,
  "millers": [
    [
      1,
      0,
      0
    ]
  ]
}

```

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```

    ]
  ],
  "elements": [
    "Al"
  ],
  "potcars": [
    "./POTCAR"
  ],
  "relax_incar" : "INCAR_metal_rlx_low",
  "scale": [
    1.0
  ],
  "skip_relax": true,
  "pert_numb": 5,
  "pert_box": 0.03,
  "pert_atom": 0.01,
  "coll_ndata": 5000,
  "_comment": "that's all"
}

```

## 5.5 dpgen init\_surf parameters

**Note:** One can load, modify, and export the input file by using our effective web-based tool [DP-GUI](#) online or hosted using the *command line interface* `dpgen gui`. All parameters below can be set in DP-GUI. By clicking “SAVE JSON”, one can download the input file.

### init\_surf\_jdata:

type: dict

argument path: `init_surf_jdata`

Generate initial data for surface systems.

#### stages:

type: `list[int]`

argument path: `init_surf_jdata/stages`

Stages for *init\_surf*.

#### elements:

type: `list[str]`

argument path: `init_surf_jdata/elements`

Atom types.

#### potcars:

type: `list[str]`, optional

argument path: `init_surf_jdata/potcars`

Path of POTCAR.

**cell\_type:**

type: str, optional

argument path: `init_surf_jdata/cell_type`

Specifying which typical structure to be generated. **Options** include fcc, hcp, bcc, sc, diamond.

**super\_cell:**

type: list[int]

argument path: `init_surf_jdata/super_cell`

Size of supercell.

**from\_poscar:**

type: bool, optional, default: False

argument path: `init_surf_jdata/from_poscar`

Deciding whether to use a given poscar as the beginning of relaxation. If it's true, keys (*cell\_type*, *latt*) will be aborted. Otherwise, these two keys are **necessary**.

**from\_poscar\_path:**

type: str, optional

argument path: `init_surf_jdata/from_poscar_path`

Path of POSCAR for VASP or STRU for ABACUS. **Necessary** if *from\_poscar* is true.

**latt:**

type: float

argument path: `init_surf_jdata/latt`

Lattice constant for single cell.

**layer\_num:**

type: int, optional

argument path: `init_surf_jdata/layer_num`

Number of atom layers constructing the slab.

**z\_min:**

type: int, optional

argument path: `init_surf_jdata/z_min`

Thickness of slab without vacuum (Angstrom). If *layer\_num* is set, *z\_min* will be ignored.

**vacuum\_max:**

type: float

argument path: `init_surf_jdata/vacuum_max`

Maximal thickness of vacuum (Angstrom).

**vacuum\_min:**

type: float, optional

argument path: `init_surf_jdata/vacuum_min`

Minimal thickness of vacuum (Angstrom). Default value is 2 times atomic radius.

**vacuum\_resol:**

type: list[float]

argument path: `init_surf_jdata/vacuum_resol`

Interval of thickness of vacuum. If size of *vacuum\_resol* is 1, the interval is fixed to its value. If size of *vacuum\_resol* is 2, the interval is *vacuum\_resol[0]* before *mid\_point*, otherwise *vacuum\_resol[1]* after *mid\_point*.

**vacuum\_numb:**

type: int, optional

argument path: `init_surf_jdata/vacuum_numb`

The total number of vacuum layers **Necessary** if *vacuum\_resol* is empty.

**mid\_point:**

type: float, optional

argument path: `init_surf_jdata/mid_point`

The mid point separating head region and tail region. **Necessary** if the size of *vacuum\_resol* is 2 or 0.

**head\_ratio:**

type: float, optional

argument path: `init_surf_jdata/head_ratio`

Ratio of vacuum layers in the nearby region with denser intervals(head region). **Necessary** if *vacuum\_resol* is empty.

**millers:**

type: list[list[int]]

argument path: `init_surf_jdata/millers`

Miller indices.

**relax\_incar:**

type: str, optional

argument path: `init_surf_jdata/relax_incar`

Path of INCAR for relaxation in VASP. **Necessary** if *stages* include 1.

**scale:**

type: list[float]

argument path: `init_surf_jdata/scale`

Scales for isotropic transforming cells.

**skip\_relax:**

type: bool

argument path: `init_surf_jdata/skip_relax`

If it's true, you may directly run stage 2 (perturb and scale) using an unrelaxed POSCAR.

**pert\_numb:**

type: int

argument path: `init_surf_jdata/pert_numb`

Number of perturbations for each scaled (key *scale*) POSCAR.

**pert\_box:**

type: float

argument path: `init_surf_jdata/pert_box`

Anisotropic Perturbation for cells (independent changes of lengths of three box vectors as well as angle among) in decimal formats. 9 elements of the 3x3 perturbation matrix will be randomly sampled from a uniform distribution (default) in the range `[-pert_box, pert_box]`. Such a perturbation matrix adds the identity matrix gives the actual transformation matrix for this perturbation operation.

**pert\_atom:**

type: float

argument path: `init_surf_jdata/pert_atom`

Perturbation of atom coordinates (Angstrom). Random perturbations are performed on three coordinates of each atom by adding values randomly sampled from a uniform distribution in the range `[-pert_atom, pert_atom]`.

**coll\_ndata:**

type: int

argument path: `init_surf_jdata/coll_ndata`

Maximal number of collected data.

## 5.6 dpgen init\_surf machine parameters

**Note:** One can load, modify, and export the input file by using our effective web-based tool [DP-GUI](#) online or hosted using the *command line interface* `dpgen gui`. All parameters below can be set in DP-GUI. By clicking “SAVE JSON”, one can download the input file.

**init\_surf\_mdata:**

type: dict

argument path: `init_surf_mdata`

machine.json file

**api\_version:**

type: str, optional, default: 1.0

argument path: `init_surf_mdata/api_version`

Please set to 1.0

**deepmd\_version:**

type: str, optional, default: 2

argument path: `init_surf_mdata/deepmd_version`

DeePMD-kit version, e.g. 2.1.3

**fp:**

type: dict

argument path: `init_surf_mdata/fp`

Parameters of command, machine, and resources for fp

**command:**

type: str  
argument path: `init_surf_mdata/fp/command`  
Command of a program.

**machine:**

type: dict  
argument path: `init_surf_mdata/fp/machine`

**batch\_type:**

type: str  
argument path: `init_surf_mdata/fp/machine/batch_type`  
The batch job system type. Option: Fugaku, Slurm, DistributedShell, SGE, PBS, Torque, SlurmJobArray, Shell, Bohrium, LSF, OpenAPI

**local\_root:**

type: str | NoneType  
argument path: `init_surf_mdata/fp/machine/local_root`  
The dir where the tasks and relating files locate. Typically the project dir.

**remote\_root:**

type: str | NoneType, optional  
argument path: `init_surf_mdata/fp/machine/remote_root`  
The dir where the tasks are executed on the remote machine. Only needed when context is not lazy-local.

**clean\_asynchronously:**

type: bool, optional, default: False  
argument path: `init_surf_mdata/fp/machine/clean_asynchronously`  
Clean the remote directory asynchronously after the job finishes.

Depending on the value of *context\_type*, different sub args are accepted.

**context\_type:**

type: str (flag key)  
argument path: `init_surf_mdata/fp/machine/context_type`



possible choices: *BohriumContext*,  
*LocalContext*, *LazyLocalContext*,  
*SSHContext*, *HDFSContext*,  
*OpenAPIContext*

The connection used to remote machine. Option: OpenAPIContext, SSHContext, BohriumContext, HDFSContext, LazyLocalContext, LocalContext

When `context_type` is set to *BohriumContext* (or its aliases *bohriumcontext*, *Bohrium*, *bohrium*, *DpCloudServerContext*, *dpcloudservercontext*, *DpCloudServer*, *dpcloudserver*, *LebesgueContext*, *lebesguecontext*, *Lebesgue*, *lebesgue*):

**remote\_profile:**

type: dict  
 argument path: `init_surf_mdata/fp/machine[BohriumContext]/remote_profile`

The information used to maintain the connection with remote machine.

**email:**

type: str, optional  
 argument path: `init_surf_mdata/fp/machine[BohriumContext]/remote_profile/email`  
 Email

**password:**

type: str, optional  
 argument path: `init_surf_mdata/fp/machine[BohriumContext]/remote_profile/password`  
 Password

**program\_id:**

type: int, alias: *project\_id*  
 argument path: `init_surf_mdata/fp/machine[BohriumContext]/remote_profile/program_id`  
 Program ID

**retry\_count:**

type: `NoneType` | int, optional,  
 default: 2  
 argument path: `init_surf_mdata/fp/machine[BohriumContext]/remote_profile/retry_count`  
 The retry count when a job is terminated

**ignore\_exit\_code:**

type: bool, optional, default: True  
argument path: init\_surf\_mdata/  
fp/machine[BohriumContext]/  
remote\_profile/  
ignore\_exit\_code

**The job state will be marked as finished if the exit code is non-zero when set to True. Otherwise,**  
the job state will be designated as terminated.

**keep\_backup:**

type: bool, optional  
argument path: init\_surf\_mdata/  
fp/machine[BohriumContext]/  
remote\_profile/keep\_backup  
  
keep download and upload zip

**input\_data:**

type: dict  
argument path: init\_surf\_mdata/  
fp/machine[BohriumContext]/  
remote\_profile/input\_data  
  
Configuration of job

When `context_type` is set to `LocalContext` (or its aliases `localcontext`, `Local`, `local`):

**remote\_profile:**

type: dict, optional  
argument path: init\_surf\_mdata/fp/  
machine[LocalContext]/  
remote\_profile

The information used to maintain the connection with remote machine. This field is empty for this context.

When `context_type` is set to `LazyLocalContext` (or its aliases `lazylocalcontext`, `LazyLocal`, `lazylocal`):

**remote\_profile:**

type: dict, optional  
argument path: init\_surf\_mdata/fp/  
machine[LazyLocalContext]/  
remote\_profile

The information used to maintain the connection with remote machine. This field is empty for this context.

When `context_type` is set to `SSHContext` (or its aliases `sshcontext`, `SSH`, `ssh`):

**remote\_profile:**

type: dict

argument path: `init_surf_mdata/fp/machine[SSHContext]/remote_profile`

The information used to maintain the connection with remote machine.

**hostname:**

type: str

argument path: `init_surf_mdata/fp/machine[SSHContext]/remote_profile/hostname`

hostname or ip of ssh connection.

**username:**

type: str

argument path: `init_surf_mdata/fp/machine[SSHContext]/remote_profile/username`

username of target linux system

**password:**

type: str, optional

argument path: `init_surf_mdata/fp/machine[SSHContext]/remote_profile/password`

(deprecated) password of linux system. Please use [SSH keys](#) instead to improve security.

**port:**

type: int, optional, default: 22

argument path: `init_surf_mdata/fp/machine[SSHContext]/remote_profile/port`

ssh connection port.

**key\_filename:**

type: str | NoneType, optional,  
default: None

argument path: `init_surf_mdata/fp/machine[SSHContext]/remote_profile/key_filename`

key filename used by ssh connection.  
If left None, find key in `~/.ssh` or use password for login

**passphrase:**

type: str | NoneType, optional,  
default: None

argument path: `init_surf_mdata/  
fp/machine[SSHContext]/  
remote_profile/passphrase`

passphrase of key used by ssh connection

**timeout:**

type: `int`, optional, default: `10`

argument path: `init_surf_mdata/  
fp/machine[SSHContext]/  
remote_profile/timeout`

timeout of ssh connection

**totp\_secret:**

type: `str | NoneType`, optional,  
default: `None`

argument path: `init_surf_mdata/  
fp/machine[SSHContext]/  
remote_profile/totp_secret`

Time-based one time password secret. It should be a base32-encoded string extracted from the 2D code.

**tar\_compress:**

type: `bool`, optional, default: `True`

argument path: `init_surf_mdata/  
fp/machine[SSHContext]/  
remote_profile/tar_compress`

The archive will be compressed in upload and download if it is `True`. If not, compression will be skipped.

**look\_for\_keys:**

type: `bool`, optional, default: `True`

argument path: `init_surf_mdata/  
fp/machine[SSHContext]/  
remote_profile/look_for_keys`

enable searching for discoverable private key files in `~/.ssh/`

When `context_type` is set to `HDFSContext` (or its aliases `hdfscontext`, `HDFS`, `hdfs`):

**remote\_profile:**

type: `dict`, optional

argument path: `init_surf_mdata/fp/  
machine[HDFSContext]/remote_profile`

The information used to maintain the connection with remote machine. This field is empty for this context.

When `context_type` is set to `OpenAPIContext` (or its aliases `openapicontext`, `OpenAPI`, `openapi`):

**remote\_profile:**

type: dict, optional  
 argument path: `init_surf_mdata/fp/machine[OpenAPIContext]/remote_profile`

The information used to maintain the connection with remote machine. This field is empty for this context.

**resources:**

type: dict  
 argument path: `init_surf_mdata/fp/resources`

**number\_node:**

type: int, optional, default: 1  
 argument path: `init_surf_mdata/fp/resources/number_node`

The number of node need for each *job*

**cpu\_per\_node:**

type: int, optional, default: 1  
 argument path: `init_surf_mdata/fp/resources/cpu_per_node`

cpu numbers of each node assigned to each job.

**gpu\_per\_node:**

type: int, optional, default: 0  
 argument path: `init_surf_mdata/fp/resources/gpu_per_node`

gpu numbers of each node assigned to each job.

**queue\_name:**

type: str, optional, default: (empty string)  
 argument path: `init_surf_mdata/fp/resources/queue_name`

The queue name of batch job scheduler system.

**group\_size:**

type: int  
 argument path: `init_surf_mdata/fp/resources/group_size`

The number of *tasks* in a *job*. 0 means infinity.

**custom\_flags:**

type: `typing.List[str]`, optional  
argument path: `init_surf_mdata/fp/  
resources/custom_flags`

The extra lines pass to job submitting script header

**strategy:**

type: `dict`, optional  
argument path: `init_surf_mdata/fp/  
resources/strategy`

strategies we use to generation job submitting scripts.

**if\_cuda\_multi\_devices:**

type: `bool`, optional, default: `False`  
argument path: `init_surf_mdata/  
fp/resources/strategy/  
if_cuda_multi_devices`

If there are multiple nvidia GPUS on the node, and we want to assign the tasks to different GPUS. If true, `dpdispatcher` will manually export environment variable `CUDA_VISIBLE_DEVICES` to different task. Usually, this option will be used with `Task.task_need_resources` variable simultaneously.

**ratio\_unfinished:**

type: `float`, optional, default: `0.0`  
argument path: `init_surf_mdata/  
fp/resources/strategy/  
ratio_unfinished`

The ratio of *tasks* that can be unfinished.

**customized\_script\_header\_template\_file:**

type: `str`, optional  
argument path: `init_surf_mdata/  
fp/resources/strategy/  
customized_script_header_template_file`

The customized template file to generate job submitting script header, which overrides the default file.

**para\_deg:**

type: `int`, optional, default: `1`  
argument path: `init_surf_mdata/fp/  
resources/para_deg`

Decide how many tasks will be run in parallel.

**source\_list:**

type: `typing.List[str]`, optional, default: `[]`

argument path: `init_surf_mdata/fp/resources/source_list`

The env file to be sourced before the command execution.

**module\_purge:**

type: `bool`, optional, default: `False`

argument path: `init_surf_mdata/fp/resources/module_purge`

Remove all modules on HPC system before module load (`module_list`)

**module\_unload\_list:**

type: `typing.List[str]`, optional, default: `[]`

argument path: `init_surf_mdata/fp/resources/module_unload_list`

The modules to be unloaded on HPC system before submitting jobs

**module\_list:**

type: `typing.List[str]`, optional, default: `[]`

argument path: `init_surf_mdata/fp/resources/module_list`

The modules to be loaded on HPC system before submitting jobs

**envs:**

type: `dict`, optional, default: `{}`

argument path: `init_surf_mdata/fp/resources/envs`

The environment variables to be exported on before submitting jobs

**prepend\_script:**

type: `typing.List[str]`, optional, default: `[]`

argument path: `init_surf_mdata/fp/resources/prepend_script`

Optional script run before jobs submitted.

**append\_script:**

type: `typing.List[str]`, optional, default: `[]`

argument path: `init_surf_mdata/fp/resources/append_script`

Optional script run after jobs submitted.

**wait\_time:**

type: `float | int`, optional, default: `0`

argument path: `init_surf_mdata/fp/resources/wait_time`

The waiting time in second after a single *task* submitted

Depending on the value of *batch\_type*, different sub args are accepted.

**batch\_type:**

type: `str` (flag key)

argument path: `init_surf_mdata/fp/resources/batch_type`

possible choices: *Bohrium*, *Fugaku*, *Slurm*, *SlurmJobArray*, *Shell*, *LSF*, *PBS*, *SGE*, *Torque*, *OpenAPI*, *DistributedShell*

The batch job system type loaded from machine/*batch\_type*.

When *batch\_type* is set to *Bohrium* (or its aliases *bohrium*, *Lebesgue*, *lebesgue*, *DpCloudServer*, *dpcloudserver*):

**kwargs:**

type: `dict`, optional

argument path: `init_surf_mdata/fp/resources[Bohrium]/kwargs`

This field is empty for this batch.

When *batch\_type* is set to *Fugaku* (or its alias *fugaku*):

**kwargs:**

type: `dict`, optional

argument path: `init_surf_mdata/fp/resources[Fugaku]/kwargs`

This field is empty for this batch.

When *batch\_type* is set to *Slurm* (or its alias *slurm*):

**kwargs:**

type: `dict`, optional



argument path: `init_surf_mdata/fp/resources[Slurm]/kwargs`

Extra arguments.

**custom\_gpu\_line:**

type: `str | NoneType`, optional,

default: `None`

argument path: `init_surf_mdata/fp/resources[Slurm]/kwargs/custom_gpu_line`

Custom GPU configuration, starting with `#SBATCH`

When `batch_type` is set to `SlurmJobArray` (or its alias `slurmjobarray`):

**kwargs:**

type: `dict`, optional

argument path: `init_surf_mdata/fp/resources[SlurmJobArray]/kwargs`

Extra arguments.

**custom\_gpu\_line:**

type: `str | NoneType`, optional,

default: `None`

argument path: `init_surf_mdata/fp/resources[SlurmJobArray]/kwargs/custom_gpu_line`

Custom GPU configuration, starting with `#SBATCH`

**slurm\_job\_size:**

type: `int`, optional, default: `1`

argument path: `init_surf_mdata/fp/resources[SlurmJobArray]/kwargs/slurm_job_size`

Number of tasks in a Slurm job

When `batch_type` is set to `Shell` (or its alias `shell`):

**kwargs:**

type: `dict`, optional

argument path: `init_surf_mdata/fp/resources[Shell]/kwargs`

This field is empty for this batch.

When `batch_type` is set to `LSF` (or its alias `lsf`):

**kwargs:**

type: `dict`

argument path: `init_surf_mdata/fp/resources[LSF]/kwargs`

Extra arguments.

**gpu\_usage:**

type: bool, optional, default: False  
argument path: `init_surf_mdata/fp/resources[LSF]/kwargs/gpu_usage`

Choosing if GPU is used in the calculation step.

**gpu\_new\_syntax:**

type: bool, optional, default: False  
argument path: `init_surf_mdata/fp/resources[LSF]/kwargs/gpu_new_syntax`

For LFS  $\geq$  10.1.0.3, new option `-gpu` for `#BSUB` could be used. If False, and old syntax would be used.

**gpu\_exclusive:**

type: bool, optional, default: True  
argument path: `init_surf_mdata/fp/resources[LSF]/kwargs/gpu_exclusive`

Only take effect when new syntax enabled. Control whether submit tasks in exclusive way for GPU.

**custom\_gpu\_line:**

type: str | NoneType, optional,  
default: None  
argument path: `init_surf_mdata/fp/resources[LSF]/kwargs/custom_gpu_line`

Custom GPU configuration, starting with `#BSUB`

When `batch_type` is set to PBS (or its alias `pbs`):

**kwargs:**

type: dict, optional  
argument path: `init_surf_mdata/fp/resources[PBS]/kwargs`

This field is empty for this batch.

When `batch_type` is set to SGE (or its alias `sge`):

**kwargs:**

type: dict, optional

argument path: `init_surf_mdata/fp/resources[SGE]/kwargs`

This field is empty for this batch.

When `batch_type` is set to Torque (or its alias `torque`):

**kwargs:**

type: dict, optional

argument path: `init_surf_mdata/fp/resources[Torque]/kwargs`

This field is empty for this batch.

When `batch_type` is set to OpenAPI (or its alias `openapi`):

**kwargs:**

type: dict, optional

argument path: `init_surf_mdata/fp/resources[OpenAPI]/kwargs`

This field is empty for this batch.

When `batch_type` is set to DistributedShell (or its alias `distributedshell`):

**kwargs:**

type: dict, optional

argument path: `init_surf_mdata/fp/resources[DistributedShell]/kwargs`

This field is empty for this batch.

**user\_forward\_files:**

type: list, optional

argument path:

`init_surf_mdata/fp/user_forward_files`

Files to be forwarded to the remote machine.

**user\_backward\_files:**

type: list, optional

argument path:

`init_surf_mdata/fp/user_backward_files`

Files to be backwarded from the remote machine.

## 5.7 init\_reaction

`dpngen init_reaction` is a workflow to initialize data for reactive systems of small gas-phase molecules. The workflow was introduced in the “Initialization” section of *Energy & Fuels*, 2021, 35 (1), 762–769.

To start the workflow, one needs a box containing reactive systems. The following packages are required for each of the step:

- Exploring: [LAMMPS](#)
- Sampling: [MDDatasetBuilder](#)
- Labeling: [Gaussian](#)

The Exploring step uses LAMMPS `pair_style reaxff` to run a short ReaxMD NVT MD simulation. In the Sampling step, molecular clusters are taken and k-means clustering algorithm is applied to remove the redundancy, which is described in *Nature Communications*, 11, 5713 (2020). The Labeling step calculates energies and forces using the Gaussian package.

An example of `reaction.json` is given below:

```

1 {
2   "type_map": [
3     "H",
4     "O"
5   ],
6   "reaxff": {
7     "data": "data.hydrogen",
8     "ff": "ffield.reax.cho",
9     "control": "lmp_control",
10    "temp": 3000,
11    "tau_t": 100,
12    "dt": 0.1,
13    "nstep": 10000,
14    "dump_freq": 100
15  },
16  "cutoff": 3.5,
17  "dataset_size": 100,
18  "qmkeywords": "b3lyp/6-31g** force Geom=PrintInputOrient"
19 }
```

For detailed parameters, see *parameters* and *machine parameters*.

The generated data can be used to continue DP-GEN concurrent learning workflow. Read *Energy & Fuels*, 2021, 35 (1), 762–769 for details.

## 5.8 dpngen init\_reaction parameters

**Note:** One can load, modify, and export the input file by using our effective web-based tool [DP-GUI](#) online or hosted using the *command line interface* `dpngen gui`. All parameters below can be set in DP-GUI. By clicking “SAVE JSON”, one can download the input file.

**init\_reaction\_jdata:**

type: dict

argument path: `init_reaction_jdata`

Generate initial data for reactive systems for small gas-phase molecules, from a ReaxFF NVT MD trajectory.

**type\_map:**

type: `list[str]`

argument path: `init_reaction_jdata/type_map`

Type map, which should match types in the initial data. e.g. [`"C"`, `"H"`, `"O"`]

**reaxff:**

type: `dict`

argument path: `init_reaction_jdata/reaxff`

Parameters for ReaxFF NVT MD.

**data:**

type: `str`

argument path: `init_reaction_jdata/reaxff/data`

Path to initial LAMMPS data file. The `atom_style` should be charge.

**ff:**

type: `str`

argument path: `init_reaction_jdata/reaxff/ff`

Path to ReaxFF force field file. Available in the `lammps/potentials` directory.

**control:**

type: `str`

argument path: `init_reaction_jdata/reaxff/control`

Path to ReaxFF control file.

**temp:**

type: `int | float`

argument path: `init_reaction_jdata/reaxff/temp`

Target Temperature for the NVT MD simulation. Unit: K.

**dt:**

type: `int | float`

argument path: `init_reaction_jdata/reaxff/dt`

Real time for every time step. Unit: fs.

**tau\_t:**

type: `int | float`

argument path: `init_reaction_jdata/reaxff/tau_t`

Time to determine how rapidly the temperature. Unit: fs.

**dump\_freq:**

type: int  
argument path: init\_reaction\_jdata/reaxff/dump\_freq  
Frequency of time steps to collect trajectory.

**nstep:**

type: int  
argument path: init\_reaction\_jdata/reaxff/nstep  
Total steps to run the ReaxFF MD simulation.

**cutoff:**

type: float  
argument path: init\_reaction\_jdata/cutoff  
Cutoff radius to take clusters from the trajectory. Note that only a complete molecule or free radical will be taken.

**dataset\_size:**

type: int  
argument path: init\_reaction\_jdata/dataset\_size  
Collected dataset size for each bond type.

**qmkeywords:**

type: str  
argument path: init\_reaction\_jdata/qmkeywords  
Gaussian keywords for first-principle calculations. e.g. force mn15/6-31g\*\*  
Geom=PrintInputOrient. Note that “force” job is necessary to collect data.  
Geom=PrintInputOrient should be used when there are more than 50 atoms in a cluster.

## 5.9 dpgen init\_reaction machine parameters

---

**Note:** One can load, modify, and export the input file by using our effective web-based tool [DP-GUI](#) online or hosted using the *command line interface* `dpgen gui`. All parameters below can be set in DP-GUI. By clicking “SAVE JSON”, one can download the input file.

---

**init\_reaction\_mdata:**

type: dict  
argument path: init\_reaction\_mdata  
machine.json file

**api\_version:**

type: str, optional, default: 1.0  
argument path: init\_reaction\_mdata/api\_version  
Please set to 1.0

**deepmd\_version:**

type: str, optional, default: 2

argument path: `init_reaction_mdata/deepmd_version`

DeePMD-kit version, e.g. 2.1.3

**reaxff:**

type: dict

argument path: `init_reaction_mdata/reaxff`

Parameters of command, machine, and resources for reaxff

**command:**

type: str

argument path: `init_reaction_mdata/reaxff/command`

Command of a program.

**machine:**

type: dict

argument path: `init_reaction_mdata/reaxff/machine`

**batch\_type:**

type: str

argument path: `init_reaction_mdata/reaxff/machine/batch_type`

The batch job system type. Option: Fugaku, Slurm, DistributedShell, SGE, PBS, Torque, SlurmJobArray, Shell, Bohrium, LSF, OpenAPI

**local\_root:**

type: str | NoneType

argument path: `init_reaction_mdata/reaxff/machine/local_root`

The dir where the tasks and relating files locate. Typically the project dir.

**remote\_root:**

type: str | NoneType, optional

argument path: `init_reaction_mdata/reaxff/machine/remote_root`

The dir where the tasks are executed on the remote machine. Only needed when context is not lazy-local.

**clean\_asynchronously:**

type: bool, optional, default: False

argument path: `init_reaction_mdata/reaxff/machine/clean_asynchronously`

Clean the remote directory asynchronously after the job finishes.

Depending on the value of *context\_type*, different sub args are accepted.

**context\_type:**

type: str (flag key)  
argument path: `init_reaction_mdata/reaxff/machine/context_type`  
possible choices: *BohriumContext*,  
*LocalContext*, *LazyLocalContext*,  
*SSHContext*, *HDFSContext*,  
*OpenAPIContext*

The connection used to remote machine. Option: *OpenAPIContext*, *SSHContext*, *BohriumContext*, *HDFSContext*, *LazyLocalContext*, *LocalContext*

When *context\_type* is set to *BohriumContext* (or its aliases *bohriumcontext*, *Bohrium*, *bohrium*, *DpCloudServerContext*, *dpcloudservercontext*, *DpCloudServer*, *dpcloudserver*, *LebesgueContext*, *lebesguecontext*, *Lebesgue*, *lebesgue*):

**remote\_profile:**

type: dict  
argument path: `init_reaction_mdata/reaxff/machine[BohriumContext]/remote_profile`

The information used to maintain the connection with remote machine.

**email:**

type: str, optional  
argument path:  
`init_reaction_mdata/reaxff/machine[BohriumContext]/remote_profile/email`  
Email

**password:**

type: str, optional  
argument path:  
`init_reaction_mdata/reaxff/machine[BohriumContext]/remote_profile/password`  
Password

**program\_id:**

type: int, alias: *project\_id*



argument path:  
 init\_reaction\_mdata/reaxff/  
 machine[BohriumContext]/  
 remote\_profile/program\_id

Program ID

#### **retry\_count:**

type: NoneType | int, optional,  
 default: 2

argument path:  
 init\_reaction\_mdata/reaxff/  
 machine[BohriumContext]/  
 remote\_profile/retry\_count

The retry count when a job is terminated

#### **ignore\_exit\_code:**

type: bool, optional, default: True

argument path:  
 init\_reaction\_mdata/reaxff/  
 machine[BohriumContext]/  
 remote\_profile/  
 ignore\_exit\_code

**The job state will be marked as finished if the exit code is non-zero when set to True. Otherwise,**  
 the job state will be designated as terminated.

#### **keep\_backup:**

type: bool, optional

argument path:  
 init\_reaction\_mdata/reaxff/  
 machine[BohriumContext]/  
 remote\_profile/keep\_backup

keep download and upload zip

#### **input\_data:**

type: dict

argument path:  
 init\_reaction\_mdata/reaxff/  
 machine[BohriumContext]/  
 remote\_profile/input\_data

Configuration of job

When `context_type` is set to LocalContext (or its aliases localcontext, Local, local):

#### **remote\_profile:**

type: dict, optional

```
argument path: init_reaction_mdata/  
reaxff/machine[LocalContext]/  
remote_profile
```

The information used to maintain the connection with remote machine. This field is empty for this context.

When `context_type` is set to `LazyLocalContext` (or its aliases `lazylocalcontext`, `LazyLocal`, `lazylocal`):

**remote\_profile:**

```
type: dict, optional  
argument path: init_reaction_mdata/  
reaxff/machine[LazyLocalContext]/  
remote_profile
```

The information used to maintain the connection with remote machine. This field is empty for this context.

When `context_type` is set to `SSHContext` (or its aliases `sshcontext`, `SSH`, `ssh`):

**remote\_profile:**

```
type: dict  
argument path:  
init_reaction_mdata/reaxff/  
machine[SSHContext]/remote_profile
```

The information used to maintain the connection with remote machine.

**hostname:**

```
type: str  
argument path:  
init_reaction_mdata/reaxff/  
machine[SSHContext]/  
remote_profile/hostname
```

hostname or ip of ssh connection.

**username:**

```
type: str  
argument path:  
init_reaction_mdata/reaxff/  
machine[SSHContext]/  
remote_profile/username
```

username of target linux system

**password:**

```
type: str, optional  
argument path:  
init_reaction_mdata/reaxff/  
machine[SSHContext]/  
remote_profile/password
```

(deprecated) password of linux system. Please use [SSH keys](#) instead to improve security.

**port:**

type: int, optional, default: 22

argument path:

init\_reaction\_mdata/reaxff/  
machine[SSHContext]/  
remote\_profile/port

ssh connection port.

**key\_filename:**

type: str | NoneType, optional,  
default: None

argument path:

init\_reaction\_mdata/reaxff/  
machine[SSHContext]/  
remote\_profile/key\_filename

key filename used by ssh connection.  
If left None, find key in ~/.ssh or use  
password for login

**passphrase:**

type: str | NoneType, optional,  
default: None

argument path:

init\_reaction\_mdata/reaxff/  
machine[SSHContext]/  
remote\_profile/passphrase

passphrase of key used by ssh connection

**timeout:**

type: int, optional, default: 10

argument path:

init\_reaction\_mdata/reaxff/  
machine[SSHContext]/  
remote\_profile/timeout

timeout of ssh connection

**totp\_secret:**

type: str | NoneType, optional,  
default: None

argument path:

init\_reaction\_mdata/reaxff/  
machine[SSHContext]/  
remote\_profile/totp\_secret

Time-based one time password secret. It should be a base32-encoded string extracted from the 2D code.

**tar\_compress:**

type: bool, optional, default: True

argument path:

init\_reaction\_mdata/reaxff/  
machine[SSHContext]/  
remote\_profile/tar\_compress

The archive will be compressed in upload and download if it is True. If not, compression will be skipped.

**look\_for\_keys:**

type: bool, optional, default: True

argument path:

init\_reaction\_mdata/reaxff/  
machine[SSHContext]/  
remote\_profile/look\_for\_keys

enable searching for discoverable private key files in ~/.ssh/

When `context_type` is set to `HDFSContext` (or its aliases `hdfscontext`, `HDFS`, `hdfs`):

**remote\_profile:**

type: dict, optional

argument path:

init\_reaction\_mdata/reaxff/  
machine[HDFSContext]/remote\_profile

The information used to maintain the connection with remote machine. This field is empty for this context.

When `context_type` is set to `OpenAPIContext` (or its aliases `openapicontext`, `OpenAPI`, `openapi`):

**remote\_profile:**

type: dict, optional

argument path: init\_reaction\_mdata/  
reaxff/machine[OpenAPIContext]/  
remote\_profile

The information used to maintain the connection with remote machine. This field is empty for this context.

**resources:**

type: dict

argument path:

init\_reaction\_mdata/reaxff/resources

**number\_node:**

type: int, optional, default: 1

argument path: `init_reaction_mdata/  
reaxff/resources/number_node`

The number of node need for each *job*

**cpu\_per\_node:**

type: int, optional, default: 1

argument path: `init_reaction_mdata/  
reaxff/resources/cpu_per_node`

cpu numbers of each node assigned to each job.

**gpu\_per\_node:**

type: int, optional, default: 0

argument path: `init_reaction_mdata/  
reaxff/resources/gpu_per_node`

gpu numbers of each node assigned to each job.

**queue\_name:**

type: str, optional, default: (empty string)

argument path: `init_reaction_mdata/  
reaxff/resources/queue_name`

The queue name of batch job scheduler system.

**group\_size:**

type: int

argument path: `init_reaction_mdata/  
reaxff/resources/group_size`

The number of *tasks* in a *job*. 0 means infinity.

**custom\_flags:**

type: typing.List[str], optional

argument path: `init_reaction_mdata/  
reaxff/resources/custom_flags`

The extra lines pass to job submitting script header

**strategy:**

type: dict, optional

argument path: `init_reaction_mdata/  
reaxff/resources/strategy`

strategies we use to generation job submitting scripts.

**if\_cuda\_multi\_devices:**

type: bool, optional, default: False

argument path:

`init_reaction_mdata/reaxff/  
resources/strategy/  
if_cuda_multi_devices`

If there are multiple nvidia GPUS on the node, and we want to assign the tasks to different GPUS. If true, dpdispatcher will manually export environment variable `CUDA_VISIBLE_DEVICES` to different task. Usually, this option will be used with `Task.task_need_resources` variable simultaneously.

**ratio\_unfinished:**

type: float, optional, default: 0.0

argument path:

`init_reaction_mdata/reaxff/  
resources/strategy/  
ratio_unfinished`

The ratio of *tasks* that can be unfinished.

**customized\_script\_header\_template\_file:**

type: str, optional

argument path:

`init_reaction_mdata/reaxff/  
resources/strategy/  
customized_script_header_template_file`

The customized template file to generate job submitting script header, which overrides the default file.

**para\_deg:**

type: int, optional, default: 1

argument path: `init_reaction_mdata/  
reaxff/resources/para_deg`

Decide how many tasks will be run in parallel.

**source\_list:**

type: `typing.List[str]`, optional, default: []

argument path: `init_reaction_mdata/  
reaxff/resources/source_list`

The env file to be sourced before the command execution.

**module\_purge:**

type: bool, optional, default: False

argument path: `init_reaction_mdata/  
reaxff/resources/module_purge`

Remove all modules on HPC system before module load (`module_list`)

**module\_unload\_list:**

type: `typing.List[str]`, optional, default: `[]`

argument path: `init_reaction_mdata/reaxff/resources/module_unload_list`

The modules to be unloaded on HPC system before submitting jobs

**module\_list:**

type: `typing.List[str]`, optional, default: `[]`

argument path: `init_reaction_mdata/reaxff/resources/module_list`

The modules to be loaded on HPC system before submitting jobs

**envs:**

type: `dict`, optional, default: `{}`

argument path: `init_reaction_mdata/reaxff/resources/envs`

The environment variables to be exported on before submitting jobs

**prepend\_script:**

type: `typing.List[str]`, optional, default: `[]`

argument path: `init_reaction_mdata/reaxff/resources/prepend_script`

Optional script run before jobs submitted.

**append\_script:**

type: `typing.List[str]`, optional, default: `[]`

argument path: `init_reaction_mdata/reaxff/resources/append_script`

Optional script run after jobs submitted.

**wait\_time:**

type: `float | int`, optional, default: `0`

argument path: `init_reaction_mdata/reaxff/resources/wait_time`

The waiting time in second after a single *task* submitted

Depending on the value of *batch\_type*, different sub args are accepted.

**batch\_type:**

type: str (flag key)  
argument path: `init_reaction_mdata/  
reaxff/resources/batch_type`  
possible choices: *Bohrium*, *Fugaku*, *Slurm*,  
*SlurmJobArray*, *Shell*, *LSF*, *PBS*, *SGE*,  
*Torque*, *OpenAPI*, *DistributedShell*

The batch job system type loaded from machine/batch\_type.

When `batch_type` is set to *Bohrium* (or its aliases *bohrium*, *Lebesgue*, *lebesgue*, *DpCloudServer*, *dpcloudserver*):

**kwargs:**

type: dict, optional  
argument path: `init_reaction_mdata/  
reaxff/resources[Bohrium]/kwargs`

This field is empty for this batch.

When `batch_type` is set to *Fugaku* (or its alias *fugaku*):

**kwargs:**

type: dict, optional  
argument path: `init_reaction_mdata/  
reaxff/resources[Fugaku]/kwargs`

This field is empty for this batch.

When `batch_type` is set to *Slurm* (or its alias *slurm*):

**kwargs:**

type: dict, optional  
argument path: `init_reaction_mdata/  
reaxff/resources[Slurm]/kwargs`

Extra arguments.

**custom\_gpu\_line:**

type: str | NoneType, optional,  
default: None  
argument path:  
`init_reaction_mdata/reaxff/  
resources[Slurm]/kwargs/  
custom_gpu_line`

Custom GPU configuration, starting with `#SBATCH`

When `batch_type` is set to *SlurmJobArray* (or its alias *slurmjobarray*):

**kwargs:**

type: dict, optional



argument path:  
 init\_reaction\_mdata/reaxff/  
 resources[SlurmJobArray]/kwargs

Extra arguments.

**custom\_gpu\_line:**

type: str | NoneType, optional,  
 default: None  
 argument path:  
 init\_reaction\_mdata/reaxff/  
 resources[SlurmJobArray]/  
 kwargs/custom\_gpu\_line  
 Custom GPU configuration, starting  
 with #SBATCH

**slurm\_job\_size:**

type: int, optional, default: 1  
 argument path:  
 init\_reaction\_mdata/reaxff/  
 resources[SlurmJobArray]/  
 kwargs/slurm\_job\_size  
 Number of tasks in a Slurm job

When `batch_type` is set to Shell (or its alias `shell`):

**kwargs:**

type: dict, optional  
 argument path: init\_reaction\_mdata/  
 reaxff/resources[Shell]/kwargs  
 This field is empty for this batch.

When `batch_type` is set to LSF (or its alias `lsf`):

**kwargs:**

type: dict  
 argument path: init\_reaction\_mdata/  
 reaxff/resources[LSF]/kwargs

Extra arguments.

**gpu\_usage:**

type: bool, optional, default: False  
 argument path:  
 init\_reaction\_mdata/reaxff/  
 resources[LSF]/kwargs/  
 gpu\_usage  
 Choosing if GPU is used in the calcu-  
 lation step.

**gpu\_new\_syntax:**

type: bool, optional, default: False

argument path:  
init\_reaction\_mdata/reaxff/  
resources[LSF]/kwargs/  
gpu\_new\_syntax

For LFS  $\geq$  10.1.0.3, new option -gpu  
for #BSUB could be used. If False,  
and old syntax would be used.

**gpu\_exclusive:**

type: bool, optional, default: True  
argument path:  
init\_reaction\_mdata/reaxff/  
resources[LSF]/kwargs/  
gpu\_exclusive

Only take effect when new syntax en-  
abled. Control whether submit tasks  
in exclusive way for GPU.

**custom\_gpu\_line:**

type: str | NoneType, optional,  
default: None  
argument path:  
init\_reaction\_mdata/reaxff/  
resources[LSF]/kwargs/  
custom\_gpu\_line

Custom GPU configuration, starting  
with #BSUB

When `batch_type` is set to PBS (or its alias pbs):

**kwargs:**

type: dict, optional  
argument path: init\_reaction\_mdata/  
reaxff/resources[PBS]/kwargs

This field is empty for this batch.

When `batch_type` is set to SGE (or its alias sge):

**kwargs:**

type: dict, optional  
argument path: init\_reaction\_mdata/  
reaxff/resources[SGE]/kwargs

This field is empty for this batch.

When `batch_type` is set to Torque (or its alias torque):

**kwargs:**

type: dict, optional  
argument path: init\_reaction\_mdata/  
reaxff/resources[Torque]/kwargs

This field is empty for this batch.

When `batch_type` is set to `OpenAPI` (or its alias `openapi`):

**kwargs:**

type: dict, optional  
 argument path: `init_reaction_mdata/reaxff/resources[OpenAPI]/kwargs`

This field is empty for this batch.

When `batch_type` is set to `DistributedShell` (or its alias `distributedshell`):

**kwargs:**

type: dict, optional  
 argument path:  
`init_reaction_mdata/reaxff/resources[DistributedShell]/kwargs`

This field is empty for this batch.

**user\_forward\_files:**

type: list, optional  
 argument path:  
`init_reaction_mdata/reaxff/user_forward_files`

Files to be forwarded to the remote machine.

**user\_backward\_files:**

type: list, optional  
 argument path:  
`init_reaction_mdata/reaxff/user_backward_files`

Files to be backwarded from the remote machine.

**build:**

type: dict  
 argument path: `init_reaction_mdata/build`  
 Parameters of command, machine, and resources for build

**command:**

type: str  
 argument path: `init_reaction_mdata/build/command`  
 Command of a program.

**machine:**

type: dict  
 argument path: `init_reaction_mdata/build/machine`

**batch\_type:**

type: str

argument path: `init_reaction_mdata/  
build/machine/batch_type`

The batch job system type. Option: Fugaku, Slurm, DistributedShell, SGE, PBS, Torque, SlurmJobArray, Shell, Bohrium, LSF, OpenAPI

#### **local\_root:**

type: `str | NoneType`

argument path: `init_reaction_mdata/  
build/machine/local_root`

The dir where the tasks and relating files locate. Typically the project dir.

#### **remote\_root:**

type: `str | NoneType`, optional

argument path: `init_reaction_mdata/  
build/machine/remote_root`

The dir where the tasks are executed on the remote machine. Only needed when context is not lazy-local.

#### **clean\_asynchronously:**

type: `bool`, optional, default: `False`

argument path: `init_reaction_mdata/  
build/machine/clean_asynchronously`

Clean the remote directory asynchronously after the job finishes.

Depending on the value of `context_type`, different sub args are accepted.

#### **context\_type:**

type: `str` (flag key)

argument path: `init_reaction_mdata/  
build/machine/context_type`

possible choices: `BohriumContext`,  
`LocalContext`, `LazyLocalContext`,  
`SSHContext`, `HDFSContext`,  
`OpenAPIContext`

The connection used to remote machine. Option: `OpenAPIContext`, `SSHContext`, `BohriumContext`, `HDFSContext`, `LazyLocalContext`, `LocalContext`

When `context_type` is set to `BohriumContext` (or its aliases `bohriumcontext`, `Bohrium`, `bohrium`, `DpCloudServerContext`, `dpcloudservercontext`, `DpCloudServer`, `dpcloudserver`, `LebesgueContext`, `lebesguecontext`, `Lebesgue`, `lebesgue`):

**remote\_profile:**

type: dict  
 argument path: init\_reaction\_mdata/  
 build/machine[BohriumContext]/  
 remote\_profile

The information used to maintain the connection with remote machine.

**email:**

type: str, optional  
 argument path:  
 init\_reaction\_mdata/build/  
 machine[BohriumContext]/  
 remote\_profile/email

Email

**password:**

type: str, optional  
 argument path:  
 init\_reaction\_mdata/build/  
 machine[BohriumContext]/  
 remote\_profile/password

Password

**program\_id:**

type: int, alias: *project\_id*  
 argument path:  
 init\_reaction\_mdata/build/  
 machine[BohriumContext]/  
 remote\_profile/program\_id

Program ID

**retry\_count:**

type: NoneType | int, optional,  
 default: 2  
 argument path:  
 init\_reaction\_mdata/build/  
 machine[BohriumContext]/  
 remote\_profile/retry\_count

The retry count when a job is terminated

**ignore\_exit\_code:**

type: bool, optional, default: True  
 argument path:  
 init\_reaction\_mdata/build/  
 machine[BohriumContext]/  
 remote\_profile/  
 ignore\_exit\_code

**The job state will be marked as finished if the exit code is non-zero when set to True. Otherwise,**  
the job state will be designated as terminated.

**keep\_backup:**

type: bool, optional  
argument path:  
init\_reaction\_mdata/build/  
machine[BohriumContext]/  
remote\_profile/keep\_backup  
keep download and upload zip

**input\_data:**

type: dict  
argument path:  
init\_reaction\_mdata/build/  
machine[BohriumContext]/  
remote\_profile/input\_data  
Configuration of job

When `context_type` is set to `LocalContext` (or its aliases `localcontext`, `Local`, `local`):

**remote\_profile:**

type: dict, optional  
argument path: init\_reaction\_mdata/  
build/machine[LocalContext]/  
remote\_profile

The information used to maintain the connection with remote machine. This field is empty for this context.

When `context_type` is set to `LazyLocalContext` (or its aliases `lazylocalcontext`, `LazyLocal`, `lazylocal`):

**remote\_profile:**

type: dict, optional  
argument path: init\_reaction\_mdata/  
build/machine[LazyLocalContext]/  
remote\_profile

The information used to maintain the connection with remote machine. This field is empty for this context.

When `context_type` is set to `SSHContext` (or its aliases `sshcontext`, `SSH`, `ssh`):

**remote\_profile:**

type: dict

argument path:  
 init\_reaction\_mdata/build/  
 machine[SSHContext]/remote\_profile

The information used to maintain the connection with remote machine.

**hostname:**

type: str  
 argument path:  
 init\_reaction\_mdata/build/  
 machine[SSHContext]/  
 remote\_profile/hostname  
 hostname or ip of ssh connection.

**username:**

type: str  
 argument path:  
 init\_reaction\_mdata/build/  
 machine[SSHContext]/  
 remote\_profile/username  
 username of target linux system

**password:**

type: str, optional  
 argument path:  
 init\_reaction\_mdata/build/  
 machine[SSHContext]/  
 remote\_profile/password  
 (deprecated) password of linux system. Please use [SSH keys](#) instead to improve security.

**port:**

type: int, optional, default: 22  
 argument path:  
 init\_reaction\_mdata/build/  
 machine[SSHContext]/  
 remote\_profile/port  
 ssh connection port.

**key\_filename:**

type: str | NoneType, optional,  
 default: None  
 argument path:  
 init\_reaction\_mdata/build/  
 machine[SSHContext]/  
 remote\_profile/key\_filename  
 key filename used by ssh connection.  
 If left None, find key in ~/.ssh or use  
 password for login

**passphrase:**

type: str | NoneType, optional,  
default: None  
argument path:  
init\_reaction\_mdata/build/  
machine[SSHContext]/  
remote\_profile/passphrase  
passphrase of key used by ssh connection

**timeout:**

type: int, optional, default: 10  
argument path:  
init\_reaction\_mdata/build/  
machine[SSHContext]/  
remote\_profile/timeout  
timeout of ssh connection

**totp\_secret:**

type: str | NoneType, optional,  
default: None  
argument path:  
init\_reaction\_mdata/build/  
machine[SSHContext]/  
remote\_profile/totp\_secret  
Time-based one time password secret. It should be a base32-encoded string extracted from the 2D code.

**tar\_compress:**

type: bool, optional, default: True  
argument path:  
init\_reaction\_mdata/build/  
machine[SSHContext]/  
remote\_profile/tar\_compress  
The archive will be compressed in upload and download if it is True. If not, compression will be skipped.

**look\_for\_keys:**

type: bool, optional, default: True  
argument path:  
init\_reaction\_mdata/build/  
machine[SSHContext]/  
remote\_profile/look\_for\_keys  
enable searching for discoverable private key files in ~/.ssh/

When `context_type` is set to `HDFSContext` (or its aliases `hdfscontext`, `HDFS`, `hdfs`):



**remote\_profile:**

type: dict, optional  
 argument path:  
 init\_reaction\_mdata/build/  
 machine[HDFSText]/remote\_profile

The information used to maintain the connection with remote machine. This field is empty for this context.

When `context_type` is set to `OpenAPIContext` (or its aliases `openapicontext`, `OpenAPI`, `openapi`):

**remote\_profile:**

type: dict, optional  
 argument path: init\_reaction\_mdata/  
 build/machine[OpenAPIContext]/  
 remote\_profile

The information used to maintain the connection with remote machine. This field is empty for this context.

**resources:**

type: dict  
 argument path:  
 init\_reaction\_mdata/build/resources

**number\_node:**

type: int, optional, default: 1  
 argument path: init\_reaction\_mdata/  
 build/resources/number\_node

The number of node need for each *job*

**cpu\_per\_node:**

type: int, optional, default: 1  
 argument path: init\_reaction\_mdata/  
 build/resources/cpu\_per\_node

cpu numbers of each node assigned to each job.

**gpu\_per\_node:**

type: int, optional, default: 0  
 argument path: init\_reaction\_mdata/  
 build/resources/gpu\_per\_node

gpu numbers of each node assigned to each job.

**queue\_name:**

type: str, optional, default: (empty string)  
 argument path: init\_reaction\_mdata/  
 build/resources/queue\_name

The queue name of batch job scheduler system.

**group\_size:**

type: int  
argument path: `init_reaction_mdata/  
build/resources/group_size`  
The number of *tasks* in a *job*. 0 means infinity.

**custom\_flags:**

type: `typing.List[str]`, optional  
argument path: `init_reaction_mdata/  
build/resources/custom_flags`  
The extra lines pass to job submitting script header

**strategy:**

type: dict, optional  
argument path: `init_reaction_mdata/  
build/resources/strategy`  
strategies we use to generation job submitting scripts.

**if\_cuda\_multi\_devices:**

type: bool, optional, default: False  
argument path:  
`init_reaction_mdata/build/  
resources/strategy/  
if_cuda_multi_devices`

If there are multiple nvidia GPUS on the node, and we want to assign the tasks to different GPUS. If true, `dpdispatcher` will manually export environment variable `CUDA_VISIBLE_DEVICES` to different task. Usually, this option will be used with `Task.task_need_resources` variable simultaneously.

**ratio\_unfinished:**

type: float, optional, default: 0.0  
argument path:  
`init_reaction_mdata/build/  
resources/strategy/  
ratio_unfinished`

The ratio of *tasks* that can be unfinished.

**customized\_script\_header\_template\_file:**

type: str, optional  
argument path:  
`init_reaction_mdata/build/`

resources/strategy/  
customized\_script\_header\_template\_file

The customized template file to generate job submitting script header, which overrides the default file.

#### **para\_deg:**

type: int, optional, default: 1  
argument path: init\_reaction\_mdata/  
build/resources/para\_deg

Decide how many tasks will be run in parallel.

#### **source\_list:**

type: typing.List[str], optional, default: []  
argument path: init\_reaction\_mdata/  
build/resources/source\_list

The env file to be sourced before the command execution.

#### **module\_purge:**

type: bool, optional, default: False  
argument path: init\_reaction\_mdata/  
build/resources/module\_purge

Remove all modules on HPC system before module load (module\_list)

#### **module\_unload\_list:**

type: typing.List[str], optional, default: []  
argument path: init\_reaction\_mdata/  
build/resources/module\_unload\_list

The modules to be unloaded on HPC system before submitting jobs

#### **module\_list:**

type: typing.List[str], optional, default: []  
argument path: init\_reaction\_mdata/  
build/resources/module\_list

The modules to be loaded on HPC system before submitting jobs

#### **envs:**

type: dict, optional, default: {}  
argument path: init\_reaction\_mdata/  
build/resources/envs

The environment variables to be exported on before submitting jobs

**prepend\_script:**

type: `typing.List[str]`, optional, default: `[]`

argument path: `init_reaction_mdata/build/resources/prepend_script`

Optional script run before jobs submitted.

**append\_script:**

type: `typing.List[str]`, optional, default: `[]`

argument path: `init_reaction_mdata/build/resources/append_script`

Optional script run after jobs submitted.

**wait\_time:**

type: `float | int`, optional, default: `0`

argument path: `init_reaction_mdata/build/resources/wait_time`

The waiting time in second after a single *task* submitted

Depending on the value of *batch\_type*, different sub args are accepted.

**batch\_type:**

type: `str` (flag key)

argument path: `init_reaction_mdata/build/resources/batch_type`

possible choices: *Bohrium*, *Fugaku*, *Slurm*, *SlurmJobArray*, *Shell*, *LSF*, *PBS*, *SGE*, *Torque*, *OpenAPI*, *DistributedShell*

The batch job system type loaded from machine/*batch\_type*.

When *batch\_type* is set to *Bohrium* (or its aliases *bohrium*, *Lebesgue*, *lebesgue*, *DpCloudServer*, *dpcloudserver*):

**kwargs:**

type: `dict`, optional

argument path: `init_reaction_mdata/build/resources[Bohrium]/kwargs`

This field is empty for this batch.

When *batch\_type* is set to *Fugaku* (or its alias *fugaku*):

**kwargs:**

type: `dict`, optional

argument path: `init_reaction_mdata/  
build/resources[Fugaku]/kwargs`

This field is empty for this batch.

When `batch_type` is set to Slurm (or its alias `slurm`):

**kwargs:**

type: dict, optional

argument path: `init_reaction_mdata/  
build/resources[Slurm]/kwargs`

Extra arguments.

**custom\_gpu\_line:**

type: str | NoneType, optional,  
default: None

argument path:  
`init_reaction_mdata/build/  
resources[Slurm]/kwargs/  
custom_gpu_line`

Custom GPU configuration, starting  
with `#SBATCH`

When `batch_type` is set to SlurmJobArray (or its alias `slurmjobarray`):

**kwargs:**

type: dict, optional

argument path:  
`init_reaction_mdata/build/  
resources[SlurmJobArray]/kwargs`

Extra arguments.

**custom\_gpu\_line:**

type: str | NoneType, optional,  
default: None

argument path:  
`init_reaction_mdata/build/  
resources[SlurmJobArray]/  
kwargs/custom_gpu_line`

Custom GPU configuration, starting  
with `#SBATCH`

**slurm\_job\_size:**

type: int, optional, default: 1

argument path:  
`init_reaction_mdata/build/  
resources[SlurmJobArray]/  
kwargs/slurm_job_size`

Number of tasks in a Slurm job

When `batch_type` is set to Shell (or its alias `shell`):

**kwargs:**

type: dict, optional  
argument path: `init_reaction_mdata/  
build/resources[Shell]/kwargs`

This field is empty for this batch.

When `batch_type` is set to LSF (or its alias `lsf`):

**kwargs:**

type: dict  
argument path: `init_reaction_mdata/  
build/resources[LSF]/kwargs`

Extra arguments.

**gpu\_usage:**

type: bool, optional, default: False  
argument path:  
`init_reaction_mdata/build/  
resources[LSF]/kwargs/  
gpu_usage`

Choosing if GPU is used in the calculation step.

**gpu\_new\_syntax:**

type: bool, optional, default: False  
argument path:  
`init_reaction_mdata/build/  
resources[LSF]/kwargs/  
gpu_new_syntax`

For LFS  $\geq$  10.1.0.3, new option `-gpu` for `#BSUB` could be used. If False, and old syntax would be used.

**gpu\_exclusive:**

type: bool, optional, default: True  
argument path:  
`init_reaction_mdata/build/  
resources[LSF]/kwargs/  
gpu_exclusive`

Only take effect when new syntax enabled. Control whether submit tasks in exclusive way for GPU.

**custom\_gpu\_line:**

type: str | NoneType, optional,  
default: None  
argument path:  
`init_reaction_mdata/build/  
resources[LSF]/kwargs/  
custom_gpu_line`

Custom GPU configuration, starting  
with #BSUB

When `batch_type` is set to PBS (or its alias `pbs`):

**kwargs:**

type: dict, optional  
argument path: `init_reaction_mdata/  
build/resources[PBS]/kwargs`

This field is empty for this batch.

When `batch_type` is set to SGE (or its alias `sge`):

**kwargs:**

type: dict, optional  
argument path: `init_reaction_mdata/  
build/resources[SGE]/kwargs`

This field is empty for this batch.

When `batch_type` is set to Torque (or its alias `torque`):

**kwargs:**

type: dict, optional  
argument path: `init_reaction_mdata/  
build/resources[Torque]/kwargs`

This field is empty for this batch.

When `batch_type` is set to OpenAPI (or its alias `openapi`):

**kwargs:**

type: dict, optional  
argument path: `init_reaction_mdata/  
build/resources[OpenAPI]/kwargs`

This field is empty for this batch.

When `batch_type` is set to DistributedShell (or its alias  
`distributedshell`):

**kwargs:**

type: dict, optional  
argument path:  
`init_reaction_mdata/build/  
resources[DistributedShell]/kwargs`

This field is empty for this batch.

**user\_forward\_files:**

type: list, optional  
argument path:  
`init_reaction_mdata/build/user_forward_files`

Files to be forwarded to the remote machine.

**user\_backward\_files:**

type: list, optional  
argument path: init\_reaction\_mdata/build/user\_backward\_files  
Files to be backwarded from the remote machine.

**fp:**

type: dict  
argument path: init\_reaction\_mdata/fp  
Parameters of command, machine, and resources for fp

**command:**

type: str  
argument path: init\_reaction\_mdata/fp/command  
Command of a program.

**machine:**

type: dict  
argument path: init\_reaction\_mdata/fp/machine

**batch\_type:**

type: str  
argument path: init\_reaction\_mdata/fp/machine/batch\_type  
The batch job system type. Option: Fugaku, Slurm, DistributedShell, SGE, PBS, Torque, SlurmJobArray, Shell, Bohrium, LSF, OpenAPI

**local\_root:**

type: str | NoneType  
argument path: init\_reaction\_mdata/fp/machine/local\_root  
The dir where the tasks and relating files locate. Typically the project dir.

**remote\_root:**

type: str | NoneType, optional  
argument path: init\_reaction\_mdata/fp/machine/remote\_root  
The dir where the tasks are executed on the remote machine. Only needed when context is not lazy-local.

**clean\_asynchronously:**

type: bool, optional, default: False



argument path: `init_reaction_mdata/fp/machine/clean_asynchronously`

Clean the remote directory asynchronously after the job finishes.

Depending on the value of `context_type`, different sub args are accepted.

#### **context\_type:**

type: str (flag key)  
 argument path: `init_reaction_mdata/fp/machine/context_type`  
 possible choices: *BohriumContext*, *LocalContext*, *LazyLocalContext*, *SSHContext*, *HDFSContext*, *OpenAPIContext*

The connection used to remote machine. Option: *OpenAPIContext*, *SSHContext*, *BohriumContext*, *HDFSContext*, *LazyLocalContext*, *LocalContext*

When `context_type` is set to *BohriumContext* (or its aliases *bohriumcontext*, *Bohrium*, *bohrium*, *DpCloudServerContext*, *dpcloudservercontext*, *DpCloudServer*, *dpcloudserver*, *LebesgueContext*, *lebesguecontext*, *Lebesgue*, *lebesgue*):

#### **remote\_profile:**

type: dict  
 argument path: `init_reaction_mdata/fp/machine[BohriumContext]/remote_profile`

The information used to maintain the connection with remote machine.

#### **email:**

type: str, optional  
 argument path:  
`init_reaction_mdata/fp/machine[BohriumContext]/remote_profile/email`

Email

#### **password:**

type: str, optional  
 argument path:  
`init_reaction_mdata/fp/machine[BohriumContext]/remote_profile/password`

Password

**program\_id:**

type: int, alias: *project\_id*  
argument path:  
init\_reaction\_mdata/fp/  
machine[BohriumContext]/  
remote\_profile/program\_id

Program ID

**retry\_count:**

type: NoneType | int, optional,  
default: 2  
argument path:  
init\_reaction\_mdata/fp/  
machine[BohriumContext]/  
remote\_profile/retry\_count

The retry count when a job is terminated

**ignore\_exit\_code:**

type: bool, optional, default: True  
argument path:  
init\_reaction\_mdata/fp/  
machine[BohriumContext]/  
remote\_profile/  
ignore\_exit\_code

**The job state will be marked as finished if the exit code is non-zero when set to True. Otherwise,** the job state will be designated as terminated.

**keep\_backup:**

type: bool, optional  
argument path:  
init\_reaction\_mdata/fp/  
machine[BohriumContext]/  
remote\_profile/keep\_backup

keep download and upload zip

**input\_data:**

type: dict  
argument path:  
init\_reaction\_mdata/fp/  
machine[BohriumContext]/  
remote\_profile/input\_data

Configuration of job

When `context_type` is set to `LocalContext` (or its aliases `localcontext`, `Local`, `local`):

**remote\_profile:**

type: dict, optional

argument path: `init_reaction_mdata/fp/  
machine[LocalContext]/  
remote_profile`

The information used to maintain the connection with remote machine. This field is empty for this context.

When `context_type` is set to `LazyLocalContext` (or its aliases `lazylocalcontext`, `LazyLocal`, `lazylocal`):

**remote\_profile:**

type: dict, optional  
argument path: `init_reaction_mdata/fp/  
machine[LazyLocalContext]/  
remote_profile`

The information used to maintain the connection with remote machine. This field is empty for this context.

When `context_type` is set to `SSHContext` (or its aliases `sshcontext`, `SSH`, `ssh`):

**remote\_profile:**

type: dict  
argument path: `init_reaction_mdata/fp/  
machine[SSHContext]/remote_profile`

The information used to maintain the connection with remote machine.

**hostname:**

type: str  
argument path:  
`init_reaction_mdata/fp/  
machine[SSHContext]/  
remote_profile/hostname`  
hostname or ip of ssh connection.

**username:**

type: str  
argument path:  
`init_reaction_mdata/fp/  
machine[SSHContext]/  
remote_profile/username`  
username of target linux system

**password:**

type: str, optional  
argument path:  
`init_reaction_mdata/fp/  
machine[SSHContext]/  
remote_profile/password`

(deprecated) password of linux system. Please use [SSH keys](#) instead to improve security.

**port:**

type: int, optional, default: 22

argument path:

init\_reaction\_mdata/fp/  
machine[SSHContext]/  
remote\_profile/port

ssh connection port.

**key\_filename:**

type: str | NoneType, optional,

default: None

argument path:

init\_reaction\_mdata/fp/  
machine[SSHContext]/  
remote\_profile/key\_filename

key filename used by ssh connection.

If left None, find key in ~/.ssh or use password for login

**passphrase:**

type: str | NoneType, optional,

default: None

argument path:

init\_reaction\_mdata/fp/  
machine[SSHContext]/  
remote\_profile/passphrase

passphrase of key used by ssh connection

**timeout:**

type: int, optional, default: 10

argument path:

init\_reaction\_mdata/fp/  
machine[SSHContext]/  
remote\_profile/timeout

timeout of ssh connection

**totp\_secret:**

type: str | NoneType, optional,

default: None

argument path:

init\_reaction\_mdata/fp/  
machine[SSHContext]/  
remote\_profile/totp\_secret

Time-based one time password secret. It should be a base32-encoded string extracted from the 2D code.

**tar\_compress:**

type: bool, optional, default: True

argument path:

init\_reaction\_mdata/fp/

machine[SSHContext]/

remote\_profile/tar\_compress

The archive will be compressed in upload and download if it is True. If not, compression will be skipped.

**look\_for\_keys:**

type: bool, optional, default: True

argument path:

init\_reaction\_mdata/fp/

machine[SSHContext]/

remote\_profile/look\_for\_keys

enable searching for discoverable private key files in ~/.ssh/

When `context_type` is set to `HDFSContext` (or its aliases `hdfscontext`, `HDFS`, `hdfs`):

**remote\_profile:**

type: dict, optional

argument path: init\_reaction\_mdata/fp/

machine[HDFSContext]/remote\_profile

The information used to maintain the connection with remote machine. This field is empty for this context.

When `context_type` is set to `OpenAPIContext` (or its aliases `openapicontext`, `OpenAPI`, `openapi`):

**remote\_profile:**

type: dict, optional

argument path: init\_reaction\_mdata/fp/

machine[OpenAPIContext]/

remote\_profile

The information used to maintain the connection with remote machine. This field is empty for this context.

**resources:**

type: dict

argument path: init\_reaction\_mdata/fp/resources

**number\_node:**

type: int, optional, default: 1

argument path: init\_reaction\_mdata/fp/

resources/number\_node

The number of node need for each *job*

**cpu\_per\_node:**

type: int, optional, default: 1  
argument path: `init_reaction_mdata/fp/resources/cpu_per_node`  
cpu numbers of each node assigned to each job.

**gpu\_per\_node:**

type: int, optional, default: 0  
argument path: `init_reaction_mdata/fp/resources/gpu_per_node`  
gpu numbers of each node assigned to each job.

**queue\_name:**

type: str, optional, default: (empty string)  
argument path: `init_reaction_mdata/fp/resources/queue_name`  
The queue name of batch job scheduler system.

**group\_size:**

type: int  
argument path: `init_reaction_mdata/fp/resources/group_size`  
The number of *tasks* in a *job*. 0 means infinity.

**custom\_flags:**

type: `typing.List[str]`, optional  
argument path: `init_reaction_mdata/fp/resources/custom_flags`  
The extra lines pass to job submitting script header

**strategy:**

type: dict, optional  
argument path: `init_reaction_mdata/fp/resources/strategy`  
strategies we use to generation job submitting scripts.

**if\_cuda\_multi\_devices:**

type: bool, optional, default: False  
argument path:  
`init_reaction_mdata/fp/resources/strategy/if_cuda_multi_devices`  
If there are multiple nvidia GPUS on the node, and we want to assign the tasks to different GPUS.If

true, `dpdispatcher` will manually export environment variable `CUDA_VISIBLE_DEVICES` to different task. Usually, this option will be used with `Task.task_need_resources` variable simultaneously.

**ratio\_unfinished:**

type: float, optional, default: 0.0  
 argument path:  
 init\_reaction\_mdata/fp/  
 resources/strategy/  
 ratio\_unfinished

The ratio of *tasks* that can be unfinished.

**customized\_script\_header\_template\_file:**

type: str, optional  
 argument path:  
 init\_reaction\_mdata/fp/  
 resources/strategy/  
 customized\_script\_header\_template\_file

The customized template file to generate job submitting script header, which overrides the default file.

**para\_deg:**

type: int, optional, default: 1  
 argument path: init\_reaction\_mdata/fp/  
 resources/para\_deg

Decide how many tasks will be run in parallel.

**source\_list:**

type: typing.List[str], optional, default: []  
 argument path: init\_reaction\_mdata/fp/  
 resources/source\_list

The env file to be sourced before the command execution.

**module\_purge:**

type: bool, optional, default: False  
 argument path: init\_reaction\_mdata/fp/  
 resources/module\_purge

Remove all modules on HPC system before module load (`module_list`)

**module\_unload\_list:**

type: typing.List[str], optional, default: []

argument path: `init_reaction_mdata/fp/resources/module_unload_list`

The modules to be unloaded on HPC system before submitting jobs

**module\_list:**

type: `typing.List[str]`, optional, default: `[]`

argument path: `init_reaction_mdata/fp/resources/module_list`

The modules to be loaded on HPC system before submitting jobs

**envs:**

type: `dict`, optional, default: `{}`

argument path: `init_reaction_mdata/fp/resources/envs`

The environment variables to be exported on before submitting jobs

**prepend\_script:**

type: `typing.List[str]`, optional, default: `[]`

argument path: `init_reaction_mdata/fp/resources/prepend_script`

Optional script run before jobs submitted.

**append\_script:**

type: `typing.List[str]`, optional, default: `[]`

argument path: `init_reaction_mdata/fp/resources/append_script`

Optional script run after jobs submitted.

**wait\_time:**

type: `float | int`, optional, default: `0`

argument path: `init_reaction_mdata/fp/resources/wait_time`

The waiting time in second after a single *task* submitted

Depending on the value of *batch\_type*, different sub args are accepted.

**batch\_type:**

type: `str` (flag key)

argument path: `init_reaction_mdata/fp/resources/batch_type`



possible choices: *Bohrium*, *Fugaku*, *Slurm*,  
*SlurmJobArray*, *Shell*, *LSF*, *PBS*, *SGE*,  
*Torque*, *OpenAPI*, *DistributedShell*

The batch job system type loaded from machine/batch\_type.

When `batch_type` is set to Bohrium (or its aliases bohrium, Lebesgue, lebesgue, DpCloudServer, dpcloudserver):

**kwargs:**

type: dict, optional  
 argument path: `init_reaction_mdata/fp/resources[Bohrium]/kwargs`

This field is empty for this batch.

When `batch_type` is set to Fugaku (or its alias fugaku):

**kwargs:**

type: dict, optional  
 argument path: `init_reaction_mdata/fp/resources[Fugaku]/kwargs`

This field is empty for this batch.

When `batch_type` is set to Slurm (or its alias slurm):

**kwargs:**

type: dict, optional  
 argument path: `init_reaction_mdata/fp/resources[Slurm]/kwargs`

Extra arguments.

**custom\_gpu\_line:**

type: str | NoneType, optional,  
 default: None  
 argument path:  
`init_reaction_mdata/fp/resources[Slurm]/kwargs/  
 custom_gpu_line`

Custom GPU configuration, starting with #SBATCH

When `batch_type` is set to SlurmJobArray (or its alias slurmjobarray):

**kwargs:**

type: dict, optional  
 argument path: `init_reaction_mdata/fp/resources[SlurmJobArray]/kwargs`

Extra arguments.

**custom\_gpu\_line:**

type: str | NoneType, optional,  
default: None

argument path:

init\_reaction\_mdata/fp/  
resources[SlurmJobArray]/  
kwargs/custom\_gpu\_line

Custom GPU configuration, starting  
with #SBATCH

**slurm\_job\_size:**

type: int, optional, default: 1

argument path:

init\_reaction\_mdata/fp/  
resources[SlurmJobArray]/  
kwargs/slurm\_job\_size

Number of tasks in a Slurm job

When `batch_type` is set to Shell (or its alias `shell`):

**kwargs:**

type: dict, optional

argument path: init\_reaction\_mdata/fp/  
resources[Shell]/kwargs

This field is empty for this batch.

When `batch_type` is set to LSF (or its alias `lsf`):

**kwargs:**

type: dict

argument path: init\_reaction\_mdata/fp/  
resources[LSF]/kwargs

Extra arguments.

**gpu\_usage:**

type: bool, optional, default: False

argument path:

init\_reaction\_mdata/fp/  
resources[LSF]/kwargs/  
gpu\_usage

Choosing if GPU is used in the calcu-  
lation step.

**gpu\_new\_syntax:**

type: bool, optional, default: False

argument path:

init\_reaction\_mdata/fp/  
resources[LSF]/kwargs/  
gpu\_new\_syntax

For LFS  $\geq$  10.1.0.3, new option -gpu for #BSUB could be used. If False, and old syntax would be used.

**gpu\_exclusive:**

type: bool, optional, default: True

argument path:

init\_reaction\_mdata/fp/

resources[LSF]/kwargs/

gpu\_exclusive

Only take effect when new syntax enabled. Control whether submit tasks in exclusive way for GPU.

**custom\_gpu\_line:**

type: str | NoneType, optional,

default: None

argument path:

init\_reaction\_mdata/fp/

resources[LSF]/kwargs/

custom\_gpu\_line

Custom GPU configuration, starting with #BSUB

When `batch_type` is set to PBS (or its alias pbs):

**kwargs:**

type: dict, optional

argument path: init\_reaction\_mdata/fp/

resources[PBS]/kwargs

This field is empty for this batch.

When `batch_type` is set to SGE (or its alias sge):

**kwargs:**

type: dict, optional

argument path: init\_reaction\_mdata/fp/

resources[SGE]/kwargs

This field is empty for this batch.

When `batch_type` is set to Torque (or its alias torque):

**kwargs:**

type: dict, optional

argument path: init\_reaction\_mdata/fp/

resources[Torque]/kwargs

This field is empty for this batch.

When `batch_type` is set to OpenAPI (or its alias openapi):

**kwargs:**

type: dict, optional

argument path: `init_reaction_mdata/fp/  
resources[OpenAPI]/kwargs`

This field is empty for this batch.

When `batch_type` is set to `DistributedShell` (or its alias `distributedshell`):

**kwargs:**

type: dict, optional  
argument path: `init_reaction_mdata/fp/  
resources[DistributedShell]/kwargs`

This field is empty for this batch.

**user\_forward\_files:**

type: list, optional  
argument path:  
`init_reaction_mdata/fp/user_forward_files`

Files to be forwarded to the remote machine.

**user\_backward\_files:**

type: list, optional  
argument path:  
`init_reaction_mdata/fp/user_backward_files`

Files to be backwarded from the remote machine.

## SIMPLIFY

## 6.1 Simplify

When you have a dataset containing lots of repeated data, this step will help you simplify your dataset. The workflow contains three stages: train, model\_devi, and fp. The train stage and the fp stage are as the same as the run step, and the model\_devi stage will calculate model deviations of the rest data that has not been confirmed accurate. Data with small model deviations will be confirmed accurate, while the program will pick data from those with large model deviations to the new dataset.

Use the following script to start the workflow:

```
dpngen simplify param.json machine.json
```

Here is an example of param.json for QM7 dataset:

```
{
  "type_map": [
    "C",
    "H",
    "N",
    "O",
    "S"
  ],
  "mass_map": [
    12.011,
    1.008,
    14.007,
    15.999,
    32.065
  ],
  "pick_data": "/scratch/jz748/simplify/qm7",
  "init_data_prefix": "",
  "init_data_sys": [],
  "sys_batch_size": [
    "auto"
  ],
  "numb_models": 4,
  "default_training_param": {
    "model": {
      "type_map": [
        "C",
```

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```

        "H",
        "N",
        "O",
        "S"
    ],
    "descriptor": {
        "type": "se_a",
        "sel": [
            7,
            16,
            3,
            3,
            1
        ],
        "rcut_smth": 1.00,
        "rcut": 6.00,
        "neuron": [
            25,
            50,
            100
        ],
        "resnet_dt": false,
        "axis_neuron": 12
    },
    "fitting_net": {
        "neuron": [
            240,
            240,
            240
        ],
        "resnet_dt": true
    }
},
"learning_rate": {
    "type": "exp",
    "start_lr": 0.001,
    "stop_lr": 5e-8,
    "decay_rate": 0.99
},
"loss": {
    "start_pref_e": 0.02,
    "limit_pref_e": 1,
    "start_pref_f": 1000,
    "limit_pref_f": 1,
    "start_pref_v": 0,
    "limit_pref_v": 0,
    "start_pref_pf": 0,
    "limit_pref_pf": 0
},
"training": {
    "set_prefix": "set",
    "numb_steps": 10000,

```

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```

        "disp_file": "lcurve.out",
        "disp_freq": 1000,
        "numb_test": 1,
        "save_freq": 1000,
        "disp_training": true,
        "time_training": true,
        "profiling": false,
        "profiling_file": "timeline.json"
    },
    "_comment": "that's all"
},
"fp_style": "gaussian",
"shuffle_poscar": false,
"fp_task_max": 1000,
"fp_task_min": 10,
"fp_pp_path": "/home/jzzeng/",
"fp_pp_files": [],
"fp_params": {
    "keywords": "mn15/6-31g** force nosymm scf(maxcyc=512)",
    "nproc": 28,
    "multiplicity": 1,
    "_comment": " that's all "
},
"init_pick_number": 100,
"iter_pick_number": 100,
"model_devi_f_trust_lo": 0.25,
"model_devi_f_trust_hi": 0.45,
"_comment": " that's all "
}

```

Here *pick\_data* is the directory to data to simplify where the program recursively detects systems System with deepmd/npz format. *init\_pick\_number* and *iter\_pick\_number* are the numbers of picked frames. *model\_devi\_f\_trust\_lo* and *model\_devi\_f\_trust\_hi* mean the range of the max deviation of atomic forces in a frame. *fp\_style* can be either gaussian or vasp currently. Other parameters are as the same as those of generator.

## 6.2 dpngen simplify parameters

**Note:** One can load, modify, and export the input file by using our effective web-based tool [DP-GUI](#) online or hosted using the *command line interface* `dpngen gui`. All parameters below can be set in DP-GUI. By clicking “SAVE JSON”, one can download the input file.

### simplify\_jdata:

type: dict

argument path: simplify\_jdata

Parameters for simplify.json, the first argument of *dpngen simplify*.

### type\_map:

type: list[str]

argument path: simplify\_jdata/type\_map

Atom types. Reminder: The elements in param.json, type.raw and data.lmp(when using lammmps) should be in the same order.

#### **mass\_map:**

type: list[float] | str, optional, default: auto

argument path: simplify\_jdata/mass\_map

Standard atomic weights (default: “auto”). if one want to use isotopes, or non-standard element names, chemical symbols, or atomic number in the type\_map list, please customize the mass\_map list instead of using “auto”.

#### **use\_ele\_temp:**

type: int, optional, default: 0

argument path: simplify\_jdata/use\_ele\_temp

Currently only support fp\_style vasp.

- 0: no electron temperature.
- 1: eletron temperature as frame parameter.
- 2: electron temperature as atom parameter.

#### **init\_data\_prefix:**

type: str, optional

argument path: simplify\_jdata/init\_data\_prefix

Prefix of initial data directories.

#### **init\_data\_sys:**

type: list[str]

argument path: simplify\_jdata/init\_data\_sys

Paths of initial data. The path can be either a system diretory containing NumPy files or an HDF5 file. You may use either absolute or relative path here. Systems will be detected recursively in the directories or the HDF5 file.

#### **sys\_format:**

type: str, optional, default: vasp/poscar

argument path: simplify\_jdata/sys\_format

Format of sys\_configs.

#### **init\_batch\_size:**

type: list[typing.Union[int, str]] | str, optional

argument path: simplify\_jdata/init\_batch\_size

Each number is the batch\_size of corresponding system for training in init\_data\_sys. One recommended rule for setting the sys\_batch\_size and init\_batch\_size is that batch\_size mutiply number of atoms ot the stucture should be larger than 32. If set to auto, batch size will be 32 divided by number of atoms. This argument will not override the mixed batch size in *default\_training\_param*.

#### **sys\_configs\_prefix:**

type: str, optional

argument path: simplify\_jdata/sys\_configs\_prefix

Prefix of sys\_configs.

#### **sys\_configs:**

type: list[list[str]]



argument path: `simplify_jdata/sys_configs`

2D list. Containing directories of structures to be explored in iterations for each system. Wildcard characters are supported here.

**sys\_batch\_size:**

type: `list[typing.Union[int, str]]`, optional

argument path: `simplify_jdata/sys_batch_size`

Each number is the `batch_size` for training of corresponding system in `sys_configs`. If set to `auto`, batch size will be 32 divided by number of atoms. This argument will not override the mixed batch size in *default\_training\_param*.

**labeled:**

type: `bool`, optional, default: `False`

argument path: `simplify_jdata/labeled`

If true, the initial data is labeled.

**pick\_data:**

type: `list[str] | str`

argument path: `simplify_jdata/pick_data`

(List of) Path to the directory with the pick data with the `deepmd/npz` or the `HDF5` file with `deepmd/hdf5` format. Systems are detected recursively.

**init\_pick\_number:**

type: `int`

argument path: `simplify_jdata/init_pick_number`

The number of initial pick data.

**iter\_pick\_number:**

type: `int`

argument path: `simplify_jdata/iter_pick_number`

The number of pick data in each iteration.

**model\_devi\_f\_trust\_lo:**

type: `float`

argument path: `simplify_jdata/model_devi_f_trust_lo`

The lower bound of forces for the selection for the model deviation.

**model\_devi\_f\_trust\_hi:**

type: `float`

argument path: `simplify_jdata/model_devi_f_trust_hi`

The higher bound of forces for the selection for the model deviation.

**model\_devi\_e\_trust\_lo:**

type: `float`, optional, default: `100000000000.0`

argument path: `simplify_jdata/model_devi_e_trust_lo`

The lower bound of energy per atom for the selection for the model deviation. Requires DeePMD-kit version  $\geq 2.2.2$ .

**model\_devi\_e\_trust\_hi:**

type: `float`, optional, default: `100000000000.0`

argument path: `simplify_jdata/model_devi_e_trust_hi`

The higher bound of energy per atom for the selection for the model deviation. Requires DeePMD-kit version  $\geq 2.2.2$ .

**true\_error\_f\_trust\_lo:**

type: float, optional, default: `10000000000.0`

argument path: `simplify_jdata/true_error_f_trust_lo`

The lower bound of forces for the selection for the true error. Requires DeePMD-kit version  $\geq 2.2.4$ .

**true\_error\_f\_trust\_hi:**

type: float, optional, default: `10000000000.0`

argument path: `simplify_jdata/true_error_f_trust_hi`

The higher bound of forces for the selection for the true error. Requires DeePMD-kit version  $\geq 2.2.4$ .

**true\_error\_e\_trust\_lo:**

type: float, optional, default: `10000000000.0`

argument path: `simplify_jdata/true_error_e_trust_lo`

The lower bound of energy per atom for the selection for the true error. Requires DeePMD-kit version  $\geq 2.2.4$ .

**true\_error\_e\_trust\_hi:**

type: float, optional, default: `10000000000.0`

argument path: `simplify_jdata/true_error_e_trust_hi`

The higher bound of energy per atom for the selection for the true error. Requires DeePMD-kit version  $\geq 2.2.4$ .

**numb\_models:**

type: int

argument path: `simplify_jdata/numb_models`

Number of models to be trained in `00.train`. 4 is recommend.

**training\_iter0\_model\_path:**

type: list[str], optional

argument path: `simplify_jdata/training_iter0_model_path`

The model used to init the first iter training. Number of element should be equal to `numb_models`.

**training\_init\_model:**

type: bool, optional

argument path: `simplify_jdata/training_init_model`

Iteration  $> 0$ , the model parameters will be initilized from the model trained at the previous iteration. Iteration  $== 0$ , the model parameters will be initialized from `training_iter0_model_path`.

**default\_training\_param:**

type: dict

argument path: `simplify_jdata/default_training_param`

Training parameters for deepmd-kit in `00.train`. You can find instructions from [DeePMD-kit documentation](#).

**dp\_train\_skip\_neighbor\_stat:**

type: bool, optional, default: False

argument path: simplify\_jdata/dp\_train\_skip\_neighbor\_stat

Append `--skip-neighbor-stat` flag to dp train.

**dp\_compress:**

type: bool, optional, default: False

argument path: simplify\_jdata/dp\_compress

Use dp compress to compress the model.

**training\_reuse\_iter:**

type: int | NoneType, optional

argument path: simplify\_jdata/training\_reuse\_iter

The minimal index of iteration that continues training models from old models of last iteration.

**training\_reuse\_old\_ratio:**

type: str | float, optional, default: auto

argument path: simplify\_jdata/training\_reuse\_old\_ratio

The probability proportion of old data during training. It can be:

- float: directly assign the probability of old data;
- *auto:f*: automatic probability, where f is the new-to-old ratio;
- *auto*: equivalent to *auto:10*.

This option is only adopted when continuing training models from old models. This option will override default parameters.

**training\_reuse\_num\_steps:**

type: int | NoneType, optional, default: None, alias:

*training\_reuse\_stop\_batch*

argument path: simplify\_jdata/training\_reuse\_num\_steps

Number of training batch. This option is only adopted when continuing training models from old models. This option will override default parameters.

**training\_reuse\_start\_lr:**

type: float | NoneType, optional, default: None

argument path: simplify\_jdata/training\_reuse\_start\_lr

The learning rate the start of the training. This option is only adopted when continuing training models from old models. This option will override default parameters.

**training\_reuse\_start\_pref\_e:**

type: int | float | NoneType, optional, default: None

argument path: simplify\_jdata/training\_reuse\_start\_pref\_e

The prefactor of energy loss at the start of the training. This option is only adopted when continuing training models from old models. This option will override default parameters.

**training\_reuse\_start\_pref\_f:**

type: int | float | NoneType, optional, default: None

argument path: `simplify_jdata/training_reuse_start_pref_f`

The prefactor of force loss at the start of the training. This option is only adopted when continuing training models from old models. This option will override default parameters.

**model\_devi\_activation\_func:**

type: `list[list[str]] | NoneType`, optional

argument path: `simplify_jdata/model_devi_activation_func`

The activation function in the model. The shape of list should be (N\_models, 2), where 2 represents the embedding and fitting network. This option will override default parameters.

**srtab\_file\_path:**

type: `str`, optional

argument path: `simplify_jdata/srtab_file_path`

The path of the table for the short-range pairwise interaction which is needed when using DP-ZBL potential

**one\_h5:**

type: `bool`, optional, default: `False`

argument path: `simplify_jdata/one_h5`

When using DeePMD-kit, all of the input data will be merged into one HDF5 file.

**training\_init\_frozen\_model:**

type: `list[str]`, optional

argument path: `simplify_jdata/training_init_frozen_model`

At iteration 0, initialize the model parameters from the given frozen models. Number of element should be equal to `numb_models`.

**training\_finetune\_model:**

type: `list[str]`, optional

argument path: `simplify_jdata/training_finetune_model`

At iteration 0, finetune the model parameters from the given frozen models. Number of element should be equal to `numb_models`.

**fp\_task\_max:**

type: `int`, optional

argument path: `simplify_jdata/fp_task_max`

Maximum of structures to be calculated in 02.fp of each iteration.

**fp\_task\_min:**

type: `int`, optional

argument path: `simplify_jdata/fp_task_min`

Minimum of structures to be calculated in 02.fp of each iteration.

**fp\_accurate\_threshold:**

type: `float`, optional

argument path: `simplify_jdata/fp_accurate_threshold`

If the accurate ratio is larger than this number, no fp calculation will be performed, i.e. `fp_task_max = 0`.

**fp\_accurate\_soft\_threshold:**

type: float, optional

argument path: `simplify_jdata/fp_accurate_soft_threshold`

If the accurate ratio is between this number and `fp_accurate_threshold`, the `fp_task_max` linearly decays to zero.

**ratio\_failed:**

type: float, optional

argument path: `simplify_jdata/ratio_failed`

Check the ratio of unsuccessfully terminated jobs. If too many FP tasks are not converged, `RuntimeError` will be raised.

Depending on the value of `fp_style`, different sub args are accepted.

**fp\_style:**

type: str (flag key), default: none

argument path: `simplify_jdata/fp_style`

possible choices: *none*, *vasp*, *gaussian*, *siesta*, *cp2k*, *abacus*, *pwmat*, *pwscf*, *custom*

Software for First Principles, if *labeled* is false.

When `fp_style` is set to *none*:

No fp.

When `fp_style` is set to *vasp*:

VASP.

**fp\_pp\_path:**

type: str

argument path: `simplify_jdata[vasp]/fp_pp_path`

Directory of psuedo-potential file to be used for 02.fp exists.

**fp\_pp\_files:**

type: list[str]

argument path: `simplify_jdata[vasp]/fp_pp_files`

Pseudo-potential file to be used for 02.fp. Note that the order of elements should correspond to the order in `type_map`.

**fp\_incar:**

type: str

argument path: `simplify_jdata[vasp]/fp_incar`

Input file for VASP. INCAR must specify KSPACING and KGAMMA.

**fp\_aniso\_kspacing:**

type: list[float], optional

argument path: `simplify_jdata[vasp]/fp_aniso_kspacing`

Set anisotropic kspacing. Usually useful for 1-D or 2-D materials. Only support VASP. If it is setting the KSPACING key in INCAR will be ignored.

**cvasp:**

type: bool, optional

argument path: `simplify_jdata[vasp]/cvasp`

If `cvasp` is true, DP-GEN will use Custodian to help control VASP calculation.

**fp\_skip\_bad\_box:**

type: str, optional

argument path: `simplify_jdata[vasp]/fp_skip_bad_box`

Skip the configurations that are obviously unreasonable before 02.fp

When `fp_style` is set to gaussian:

Gaussian. The command should be set as `g16 < input`.

**use\_clusters:**

type: bool, optional, default: False

argument path: `simplify_jdata[gaussian]/use_clusters`

If set to true, clusters will be taken instead of the whole system.

**cluster\_cutoff:**

type: float, optional

argument path: `simplify_jdata[gaussian]/cluster_cutoff`

The soft cutoff radius of clusters if `use_clusters` is set to true. Molecules will be taken as whole even if part of atoms is out of the cluster. Use `cluster_cutoff_hard` to only take atoms within the hard cutoff radius.

**cluster\_cutoff\_hard:**

type: float, optional

argument path: `simplify_jdata[gaussian]/cluster_cutoff_hard`

The hard cutoff radius of clusters if `use_clusters` is set to true. Outside the hard cutoff radius, atoms will not be taken even if they are in a molecule where some atoms are within the cutoff radius.

**cluster\_minify:**

type: bool, optional, default: False

argument path: `simplify_jdata[gaussian]/cluster_minify`

If enabled, when an atom within the soft cutoff radius connects a single bond with a non-hydrogen atom out of the soft cutoff radius, the outer atom will be replaced by a hydrogen atom. When the outer atom is a hydrogen atom, the outer atom will be kept. In this case, other atoms out of the soft cutoff radius will be removed.

**fp\_params:**

type: dict

argument path: `simplify_jdata[gaussian]/fp_params`

Parameters for Gaussian calculation.

**keywords:**

type: `list[str] | str`

argument path:

`simplify_jdata[gaussian]/fp_params/keywords`

Keywords for Gaussian input, e.g. force b3lyp/6-31g\*\*. If a list, run multiple steps.

#### **multiplicity:**

type: str | int, optional, default: auto  
 argument path: simplify\_jdata[gaussian]/  
 fp\_params/multiplicity

Spin multiplicity for Gaussian input. If *auto*, multiplicity will be detected automatically, with the following rules: when fragment\_guesses=True, multiplicity will +1 for each radical, and +2 for each oxygen molecule; when fragment\_guesses=False, multiplicity will be 1 or 2, but +2 for each oxygen molecule.

#### **nproc:**

type: int  
 argument path:  
 simplify\_jdata[gaussian]/fp\_params/nproc

The number of processors for Gaussian input.

#### **charge:**

type: int, optional, default: 0  
 argument path:  
 simplify\_jdata[gaussian]/fp\_params/charge

Molecule charge. Only used when charge is not provided by the system.

#### **fragment\_guesses:**

type: bool, optional, default: False  
 argument path: simplify\_jdata[gaussian]/  
 fp\_params/fragment\_guesses

Initial guess generated from fragment guesses. If True, *multiplicity* should be *auto*.

#### **basis\_set:**

type: str, optional  
 argument path:  
 simplify\_jdata[gaussian]/fp\_params/basis\_set

Custom basis set.

#### **keywords\_high\_multiplicity:**

type: str, optional  
 argument path: simplify\_jdata[gaussian]/  
 fp\_params/keywords\_high\_multiplicity

Keywords for points with multiple radicals. *multiplicity* should be *auto*. If not set, fallback to normal keywords.

When `fp_style` is set to `siesta`:

**use\_clusters:**

type: bool, optional

argument path: `simplify_jdata[siesta]/use_clusters`

If set to true, clusters will be taken instead of the whole system. This option does not work with DeePMD-kit 0.x.

**cluster\_cutoff:**

type: float, optional

argument path: `simplify_jdata[siesta]/cluster_cutoff`

The cutoff radius of clusters if `use_clusters` is set to true.

**fp\_params:**

type: dict

argument path: `simplify_jdata[siesta]/fp_params`

Parameters for siesta calculation.

**ecut:**

type: int

argument path:

`simplify_jdata[siesta]/fp_params/ecut`

Define the plane wave cutoff for grid.

**ediff:**

type: float

argument path:

`simplify_jdata[siesta]/fp_params/ediff`

Tolerance of Density Matrix.

**kspacing:**

type: float

argument path:

`simplify_jdata[siesta]/fp_params/kspacing`

Sample factor in Brillouin zones.

**mixingWeight:**

type: float

argument path:

`simplify_jdata[siesta]/fp_params/mixingWeight`

Proportion a of output Density Matrix to be used for the input Density Matrix of next SCF cycle (linear mixing).

**NumberPulay:**

type: int

argument path:

`simplify_jdata[siesta]/fp_params/NumberPulay`

Controls the Pulay convergence accelerator.



**fp\_pp\_path:**

type: str

argument path: `simplify_jdata[siesta]/fp_pp_path`

Directory of psuedo-potential or numerical orbital files to be used for 02.fp exists.

**fp\_pp\_files:**

type: list[str]

argument path: `simplify_jdata[siesta]/fp_pp_files`Psuedo-potential file to be used for 02.fp. Note that the order of elements should correspond to the order in `type_map`.When `fp_style` is set to `cp2k`:**user\_fp\_params:**type: dict, optional, alias: `fp_params`argument path: `simplify_jdata[cp2k]/user_fp_params`Parameters for cp2k calculation. find detail in [manual.cp2k.org](http://manual.cp2k.org). only the kind section must be set before use. we assume that you have basic knowledge for cp2k input.**external\_input\_path:**

type: str, optional

argument path: `simplify_jdata[cp2k]/external_input_path`Conflict with key: `user_fp_params`. enable the template input provided by user. some rules should be followed, read the following text in detail:

1. One must present a KEYWORD ABC in the section CELL so that the DP-GEN can replace the cell on-the-fly.
2. One need to add these lines under FORCE\_EVAL section to print forces and stresses:

```
STRESS_TENSOR ANALYTICAL
&PRINT
  &FORCES ON
  &END FORCES
  &STRESS_TENSOR ON
  &END STRESS_TENSOR
&END PRINT
```

When `fp_style` is set to `abacus`:**fp\_pp\_path:**

type: str

argument path: `simplify_jdata[abacus]/fp_pp_path`

Directory of psuedo-potential or numerical orbital files to be used for 02.fp exists.

**fp\_pp\_files:**

type: list[str]

argument path: `simplify_jdata[abacus]/fp_pp_files`Psuedo-potential file to be used for 02.fp. Note that the order of elements should correspond to the order in `type_map`.

**fp\_orb\_files:**

type: list[str], optional

argument path: simplify\_jdata[abacus]/fp\_orb\_files

numerical orbital file to be used for 02.fp when using LCAO basis. Note that the order of elements should correspond to the order in type\_map.

**fp\_incar:**

type: str, optional

argument path: simplify\_jdata[abacus]/fp\_incar

Input file for ABACUS. This is optional but the priority is lower than user\_fp\_params, and you should not set user\_fp\_params if you want to use fp\_incar.

**fp\_kpt\_file:**

type: str, optional

argument path: simplify\_jdata[abacus]/fp\_kpt\_file

KPT file for ABACUS. If the “kspacing” or “gamma\_only=1” is defined in INPUT or “k\_points” is defined, fp\_kpt\_file will be ignored.

**fp\_dpks\_descriptor:**

type: str, optional

argument path: simplify\_jdata[abacus]/fp\_dpks\_descriptor

DeePKS descriptor file name. The file should be in pseudopotential directory.

**user\_fp\_params:**

type: dict, optional

argument path: simplify\_jdata[abacus]/user\_fp\_params

Set the key and value of INPUT.

**k\_points:**

type: list[int], optional

argument path: simplify\_jdata[abacus]/k\_points

Monkhorst-Pack k-grids setting for generating KPT file of ABACUS, such as: [1,1,1,0,0,0]. NB: if “kspacing” or “gamma\_only=1” is defined in INPUT, k\_points will be ignored.

When `fp_style` is set to `pwmat`:

TODO: add doc

When `fp_style` is set to `pwscf`:

`pwscf` (Quantum Espresso).

**fp\_pp\_path:**

type: str

argument path: simplify\_jdata[pwscf]/fp\_pp\_path

Directory of psuedo-potential file to be used for 02.fp exists.

**fp\_pp\_files:**

type: list[str]

argument path: `simplify_jdata[pwscf]/fp_pp_files`

Pseudo-potential file to be used for 02.fp. Note that the order of elements should correspond to the order in `type_map`.

#### **fp\_params:**

type: dict, optional

argument path: `simplify_jdata[pwscf]/fp_params`

Parameters for pwscf calculation. It has lower priority than `user_fp_params`.

#### **ecut:**

type: float

argument path:

`simplify_jdata[pwscf]/fp_params/ecut`

`ecutwfc` in pwscf.

#### **ediff:**

type: float

argument path:

`simplify_jdata[pwscf]/fp_params/ediff`

`conv_thr` and `ts_vdw_econv_thr` in pwscf.

#### **smearing:**

type: str

argument path:

`simplify_jdata[pwscf]/fp_params/smearing`

`smearing` in pwscf.

#### **sigma:**

type: float

argument path:

`simplify_jdata[pwscf]/fp_params/sigma`

`degauss` in pwscf.

#### **kspacing:**

type: float

argument path:

`simplify_jdata[pwscf]/fp_params/kspacing`

The spacing between kpoints. Helps to determin KPOINTS in pwscf.

#### **user\_fp\_params:**

type: dict, optional

argument path: `simplify_jdata[pwscf]/user_fp_params`

Parameters for pwscf calculation. Find details at [https://www.quantum-espresso.org/Doc/INPUT\\_PW.html](https://www.quantum-espresso.org/Doc/INPUT_PW.html). When `user_fp_params` is set, the settings in `fp_params` will be ignored. If one wants to use `user_fp_params`, `kspacing` must be set in `user_fp_params`. `kspacing` is the spacing between kpoints, and helps to determin KPOINTS in pwscf.

When `fp_style` is set to `custom`:

Custom FP code. You need to provide the input and output file format and name. The command argument in the machine file should be the script to run custom FP codes. The extra forward and backward files can be defined in the machine file.

**fp\_params:**

type: dict  
argument path: `simplify_jdata[custom]/fp_params`

Parameters for FP calculation.

**input\_fmt:**

type: str  
argument path:  
`simplify_jdata[custom]/fp_params/input_fmt`  
Input dpdata format of the custom FP code. Such format should only need the first argument as the file name.

**input\_fn:**

type: str  
argument path:  
`simplify_jdata[custom]/fp_params/input_fn`  
Input file name of the custom FP code.

**output\_fmt:**

type: str  
argument path:  
`simplify_jdata[custom]/fp_params/output_fmt`  
Output dpdata format of the custom FP code. Such format should only need the first argument as the file name.

**output\_fn:**

type: str  
argument path:  
`simplify_jdata[custom]/fp_params/output_fn`  
Output file name of the custom FP code.

## 6.3 dpgen simplify machine parameters

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**Note:** One can load, modify, and export the input file by using our effective web-based tool [DP-GUI](#) online or hosted using the *command line interface* `dpgen gui`. All parameters below can be set in DP-GUI. By clicking “SAVE JSON”, one can download the input file.

---

**simplify\_mdata:**

type: dict  
argument path: `simplify_mdata`  
machine.json file

**api\_version:**

type: str, optional, default: 1.0  
 argument path: simplify\_mdata/api\_version  
 Please set to 1.0

**deepmd\_version:**

type: str, optional, default: 2  
 argument path: simplify\_mdata/deepmd\_version  
 DeePMD-kit version, e.g. 2.1.3

**train:**

type: dict  
 argument path: simplify\_mdata/train  
 Parameters of command, machine, and resources for train

**command:**

type: str  
 argument path: simplify\_mdata/train/command  
 Command of a program.

**machine:**

type: dict  
 argument path: simplify\_mdata/train/machine

**batch\_type:**

type: str  
 argument path: simplify\_mdata/train/machine/batch\_type  
 The batch job system type. Option: Fugaku, Slurm, DistributedShell, SGE, PBS, Torque, SlurmJobArray, Shell, Bohrium, LSF, OpenAPI

**local\_root:**

type: str | NoneType  
 argument path: simplify\_mdata/train/machine/local\_root  
 The dir where the tasks and relating files locate. Typically the project dir.

**remote\_root:**

type: str | NoneType, optional  
 argument path: simplify\_mdata/train/machine/remote\_root  
 The dir where the tasks are executed on the remote machine. Only needed when context is not lazy-local.

**clean\_asynchronously:**

type: bool, optional, default: False  
argument path: `simplify_mdata/train/  
machine/clean_asynchronously`

Clean the remote directory asynchronously after the job finishes.

Depending on the value of `context_type`, different sub args are accepted.

**context\_type:**

type: str (flag key)  
argument path: `simplify_mdata/train/  
machine/context_type`  
possible choices: `BohriumContext`,  
`LocalContext`, `LazyLocalContext`,  
`SSHContext`, `HDFSContext`,  
`OpenAPIContext`

The connection used to remote machine. Option: `OpenAPIContext`, `SSHContext`, `BohriumContext`, `HDFSContext`, `LazyLocalContext`, `LocalContext`

When `context_type` is set to `BohriumContext` (or its aliases `bohriumcontext`, `Bohrium`, `bohrium`, `DpCloudServerContext`, `dpcloudservercontext`, `DpCloudServer`, `dpcloudserver`, `LebesgueContext`, `lebesguecontext`, `Lebesgue`, `lebesgue`):

**remote\_profile:**

type: dict  
argument path: `simplify_mdata/train/  
machine[BohriumContext]/  
remote_profile`

The information used to maintain the connection with remote machine.

**email:**

type: str, optional  
argument path:  
`simplify_mdata/train/  
machine[BohriumContext]/  
remote_profile/email`

Email

**password:**

type: str, optional  
argument path:  
`simplify_mdata/train/  
machine[BohriumContext]/  
remote_profile/password`

Password

**program\_id:**

type: int, alias: *project\_id*  
 argument path:  
 simplify\_mdata/train/  
 machine[BohriumContext]/  
 remote\_profile/program\_id

Program ID

**retry\_count:**

type: NoneType | int, optional,  
 default: 2  
 argument path:  
 simplify\_mdata/train/  
 machine[BohriumContext]/  
 remote\_profile/retry\_count

The retry count when a job is terminated

**ignore\_exit\_code:**

type: bool, optional, default: True  
 argument path:  
 simplify\_mdata/train/  
 machine[BohriumContext]/  
 remote\_profile/  
 ignore\_exit\_code

**The job state will be marked as finished if the exit code is non-zero when set to True. Otherwise,**  
 the job state will be designated as terminated.

**keep\_backup:**

type: bool, optional  
 argument path:  
 simplify\_mdata/train/  
 machine[BohriumContext]/  
 remote\_profile/keep\_backup

keep download and upload zip

**input\_data:**

type: dict  
 argument path:  
 simplify\_mdata/train/  
 machine[BohriumContext]/  
 remote\_profile/input\_data

Configuration of job

When `context_type` is set to `LocalContext` (or its aliases `localcontext`, `Local`, `local`):

**remote\_profile:**

type: dict, optional  
argument path: simplify\_mdata/train/  
machine[LocalContext]/  
remote\_profile

The information used to maintain the connection with remote machine. This field is empty for this context.

When `context_type` is set to `LazyLocalContext` (or its aliases `lazylocalcontext`, `LazyLocal`, `lazylocal`):

**remote\_profile:**

type: dict, optional  
argument path: simplify\_mdata/train/  
machine[LazyLocalContext]/  
remote\_profile

The information used to maintain the connection with remote machine. This field is empty for this context.

When `context_type` is set to `SSHContext` (or its aliases `sshcontext`, `SSH`, `ssh`):

**remote\_profile:**

type: dict  
argument path: simplify\_mdata/train/  
machine[SSHContext]/remote\_profile

The information used to maintain the connection with remote machine.

**hostname:**

type: str  
argument path: simplify\_mdata/  
train/machine[SSHContext]/  
remote\_profile/hostname

hostname or ip of ssh connection.

**username:**

type: str  
argument path: simplify\_mdata/  
train/machine[SSHContext]/  
remote\_profile/username

username of target linux system

**password:**

type: str, optional  
argument path: simplify\_mdata/  
train/machine[SSHContext]/  
remote\_profile/password



(deprecated) password of linux system. Please use [SSH keys](#) instead to improve security.

**port:**

type: int, optional, default: 22  
 argument path: `simplify_mdata/  
 train/machine[SSHContext]/  
 remote_profile/port`  
 ssh connection port.

**key\_filename:**

type: str | NoneType, optional,  
 default: None  
 argument path: `simplify_mdata/  
 train/machine[SSHContext]/  
 remote_profile/key_filename`  
 key filename used by ssh connection.  
 If left None, find key in `~/.ssh` or use  
 password for login

**passphrase:**

type: str | NoneType, optional,  
 default: None  
 argument path: `simplify_mdata/  
 train/machine[SSHContext]/  
 remote_profile/passphrase`  
 passphrase of key used by ssh connection

**timeout:**

type: int, optional, default: 10  
 argument path: `simplify_mdata/  
 train/machine[SSHContext]/  
 remote_profile/timeout`  
 timeout of ssh connection

**totp\_secret:**

type: str | NoneType, optional,  
 default: None  
 argument path: `simplify_mdata/  
 train/machine[SSHContext]/  
 remote_profile/totp_secret`  
 Time-based one time password secret. It should be a base32-encoded string extracted from the 2D code.

**tar\_compress:**

type: bool, optional, default: True  
 argument path: `simplify_mdata/  
 train/machine[SSHContext]/  
 remote_profile/tar_compress`

The archive will be compressed in upload and download if it is True. If not, compression will be skipped.

**look\_for\_keys:**

type: bool, optional, default: True  
argument path: `simplify_mdata/train/machine[SSHContext]/remote_profile/look_for_keys`  
enable searching for discoverable private key files in `~/.ssh/`

When `context_type` is set to `HDFSContext` (or its aliases `hdfscontext`, `HDFS`, `hdfs`):

**remote\_profile:**

type: dict, optional  
argument path: `simplify_mdata/train/machine[HDFSContext]/remote_profile`  
The information used to maintain the connection with remote machine. This field is empty for this context.

When `context_type` is set to `OpenAPIContext` (or its aliases `openapicontext`, `OpenAPI`, `openapi`):

**remote\_profile:**

type: dict, optional  
argument path: `simplify_mdata/train/machine[OpenAPIContext]/remote_profile`  
The information used to maintain the connection with remote machine. This field is empty for this context.

**resources:**

type: dict  
argument path: `simplify_mdata/train/resources`

**number\_node:**

type: int, optional, default: 1  
argument path: `simplify_mdata/train/resources/number_node`  
The number of node need for each *job*

**cpu\_per\_node:**

type: int, optional, default: 1  
argument path: `simplify_mdata/train/resources/cpu_per_node`  
cpu numbers of each node assigned to each job.

**gpu\_per\_node:**

type: int, optional, default: 0

argument path: `simplify_mdata/train/resources/gpu_per_node`

gpu numbers of each node assigned to each job.

**queue\_name:**

type: str, optional, default: (empty string)

argument path: `simplify_mdata/train/resources/queue_name`

The queue name of batch job scheduler system.

**group\_size:**

type: int

argument path: `simplify_mdata/train/resources/group_size`

The number of *tasks* in a *job*. 0 means infinity.

**custom\_flags:**

type: `typing.List[str]`, optional

argument path: `simplify_mdata/train/resources/custom_flags`

The extra lines pass to job submitting script header

**strategy:**

type: dict, optional

argument path: `simplify_mdata/train/resources/strategy`

strategies we use to generation job submitting scripts.

**if\_cuda\_multi\_devices:**

type: bool, optional, default: False

argument path: `simplify_mdata/train/resources/strategy/if_cuda_multi_devices`

If there are multiple nvidia GPUS on the node, and we want to assign the tasks to different GPUS. If true, `dpdispatcher` will manually export environment variable `CUDA_VISIBLE_DEVICES` to different task. Usually, this option will be used with `Task.task_need_resources` variable simultaneously.

**ratio\_unfinished:**

type: float, optional, default: 0.0  
argument path: simplify\_mdata/  
train/resources/strategy/  
ratio\_unfinished

The ratio of *tasks* that can be unfinished.

**customized\_script\_header\_template\_file:**

type: str, optional  
argument path: simplify\_mdata/  
train/resources/strategy/  
customized\_script\_header\_template\_file

The customized template file to generate job submitting script header, which overrides the default file.

**para\_deg:**

type: int, optional, default: 1  
argument path: simplify\_mdata/train/  
resources/para\_deg

Decide how many tasks will be run in parallel.

**source\_list:**

type: typing.List[str], optional, default: []  
argument path: simplify\_mdata/train/  
resources/source\_list

The env file to be sourced before the command execution.

**module\_purge:**

type: bool, optional, default: False  
argument path: simplify\_mdata/train/  
resources/module\_purge

Remove all modules on HPC system before module load (module\_list)

**module\_unload\_list:**

type: typing.List[str], optional, default: []  
argument path: simplify\_mdata/train/  
resources/module\_unload\_list

The modules to be unloaded on HPC system before submitting jobs

**module\_list:**

type: typing.List[str], optional, default: []

argument path: `simplify_mdata/train/resources/module_list`

The modules to be loaded on HPC system before submitting jobs

#### **envs:**

type: dict, optional, default: {}

argument path: `simplify_mdata/train/resources/envs`

The environment variables to be exported on before submitting jobs

#### **prepend\_script:**

type: `typing.List[str]`, optional, default: []

argument path: `simplify_mdata/train/resources/prepend_script`

Optional script run before jobs submitted.

#### **append\_script:**

type: `typing.List[str]`, optional, default: []

argument path: `simplify_mdata/train/resources/append_script`

Optional script run after jobs submitted.

#### **wait\_time:**

type: `float | int`, optional, default: 0

argument path: `simplify_mdata/train/resources/wait_time`

The waiting time in second after a single *task* submitted

Depending on the value of *batch\_type*, different sub args are accepted.

#### **batch\_type:**

type: str (flag key)

argument path: `simplify_mdata/train/resources/batch_type`

possible choices: *Bohrium*, *Fugaku*, *Slurm*, *SlurmJobArray*, *Shell*, *LSF*, *PBS*, *SGE*, *Torque*, *OpenAPI*, *DistributedShell*

The batch job system type loaded from machine/*batch\_type*.

When *batch\_type* is set to *Bohrium* (or its aliases *bohrium*, *Lebesgue*, *lebesgue*, *DpCloudServer*, *dpcloudserver*):

**kwargs:**

type: dict, optional  
argument path: `simplify_mdata/train/resources[Bohrium]/kwargs`

This field is empty for this batch.

When `batch_type` is set to Fugaku (or its alias `fugaku`):

**kwargs:**

type: dict, optional  
argument path: `simplify_mdata/train/resources[Fugaku]/kwargs`

This field is empty for this batch.

When `batch_type` is set to Slurm (or its alias `slurm`):

**kwargs:**

type: dict, optional  
argument path: `simplify_mdata/train/resources[Slurm]/kwargs`

Extra arguments.

**custom\_gpu\_line:**

type: str | NoneType, optional,  
default: None  
argument path: `simplify_mdata/train/resources[Slurm]/kwargs/custom_gpu_line`  
  
Custom GPU configuration, starting  
with `#SBATCH`

When `batch_type` is set to SlurmJobArray (or its alias `slurmjobarray`):

**kwargs:**

type: dict, optional  
argument path: `simplify_mdata/train/resources[SlurmJobArray]/kwargs`

Extra arguments.

**custom\_gpu\_line:**

type: str | NoneType, optional,  
default: None  
argument path:  
`simplify_mdata/train/resources[SlurmJobArray]/kwargs/custom_gpu_line`  
  
Custom GPU configuration, starting  
with `#SBATCH`

**slurm\_job\_size:**

type: int, optional, default: 1  
 argument path:  
 simplify\_mdata/train/  
 resources[SlurmJobArray]/  
 kwargs/slurm\_job\_size

Number of tasks in a Slurm job

When `batch_type` is set to Shell (or its alias `shell`):

**kwargs:**

type: dict, optional  
 argument path: simplify\_mdata/train/  
 resources[Shell]/kwargs

This field is empty for this batch.

When `batch_type` is set to LSF (or its alias `lsf`):

**kwargs:**

type: dict  
 argument path: simplify\_mdata/train/  
 resources[LSF]/kwargs

Extra arguments.

**gpu\_usage:**

type: bool, optional, default: False  
 argument path: simplify\_mdata/  
 train/resources[LSF]/kwargs/  
 gpu\_usage

Choosing if GPU is used in the calculation step.

**gpu\_new\_syntax:**

type: bool, optional, default: False  
 argument path: simplify\_mdata/  
 train/resources[LSF]/kwargs/  
 gpu\_new\_syntax

For LFS >= 10.1.0.3, new option -gpu for #BSUB could be used. If False, and old syntax would be used.

**gpu\_exclusive:**

type: bool, optional, default: True  
 argument path: simplify\_mdata/  
 train/resources[LSF]/kwargs/  
 gpu\_exclusive

Only take effect when new syntax enabled. Control whether submit tasks in exclusive way for GPU.

**custom\_gpu\_line:**

type: str | NoneType, optional,  
default: None  
argument path: simplify\_mdata/  
train/resources[LSF]/kwargs/  
custom\_gpu\_line  
Custom GPU configuration, starting  
with #BSUB

When `batch_type` is set to PBS (or its alias pbs):

**kwargs:**

type: dict, optional  
argument path: simplify\_mdata/train/  
resources[PBS]/kwargs  
This field is empty for this batch.

When `batch_type` is set to SGE (or its alias sge):

**kwargs:**

type: dict, optional  
argument path: simplify\_mdata/train/  
resources[SGE]/kwargs  
This field is empty for this batch.

When `batch_type` is set to Torque (or its alias torque):

**kwargs:**

type: dict, optional  
argument path: simplify\_mdata/train/  
resources[Torque]/kwargs  
This field is empty for this batch.

When `batch_type` is set to OpenAPI (or its alias openapi):

**kwargs:**

type: dict, optional  
argument path: simplify\_mdata/train/  
resources[OpenAPI]/kwargs  
This field is empty for this batch.

When `batch_type` is set to DistributedShell (or its alias distributedshell):

**kwargs:**

type: dict, optional  
argument path: simplify\_mdata/train/  
resources[DistributedShell]/kwargs  
This field is empty for this batch.



**user\_forward\_files:**

type: list, optional  
 argument path: `simplify_mdata/train/user_forward_files`  
 Files to be forwarded to the remote machine.

**user\_backward\_files:**

type: list, optional  
 argument path: `simplify_mdata/train/user_backward_files`  
 Files to be backwarded from the remote machine.

**model\_devi:**

type: dict  
 argument path: `simplify_mdata/model_devi`  
 Parameters of command, machine, and resources for model\_devi

**command:**

type: str  
 argument path: `simplify_mdata/model_devi/command`  
 Command of a program.

**machine:**

type: dict  
 argument path: `simplify_mdata/model_devi/machine`

**batch\_type:**

type: str  
 argument path: `simplify_mdata/model_devi/machine/batch_type`  
 The batch job system type. Option: Fugaku, Slurm, DistributedShell, SGE, PBS, Torque, SlurmJobArray, Shell, Bohrium, LSF, OpenAPI

**local\_root:**

type: str | NoneType  
 argument path: `simplify_mdata/model_devi/machine/local_root`  
 The dir where the tasks and relating files locate. Typically the project dir.

**remote\_root:**

type: str | NoneType, optional  
 argument path: `simplify_mdata/model_devi/machine/remote_root`

The dir where the tasks are executed on the remote machine. Only needed when context is not lazy-local.

**clean\_asynchronously:**

type: bool, optional, default: False

argument path:

simplify\_mdata/model\_devi/machine/  
clean\_asynchronously

Clean the remote directory asynchronously after the job finishes.

Depending on the value of *context\_type*, different sub args are accepted.

**context\_type:**

type: str (flag key)

argument path: simplify\_mdata/  
model\_devi/machine/context\_type

possible choices: *BohriumContext*,  
*LocalContext*, *LazyLocalContext*,  
*SSHContext*, *HDFSContext*,  
*OpenAPIContext*

The connection used to remote machine. Option: OpenAPIContext, SSHContext, BohriumContext, HDFSContext, LazyLocalContext, LocalContext

When *context\_type* is set to *BohriumContext* (or its aliases *bohriumcontext*, *Bohrium*, *bohrium*, *DpCloudServerContext*, *dpcloudservercontext*, *DpCloudServer*, *dpcloudserver*, *LebesgueContext*, *lebesguecontext*, *Lebesgue*, *lebesgue*):

**remote\_profile:**

type: dict

argument path: simplify\_mdata/  
model\_devi/machine[BohriumContext]/  
remote\_profile

The information used to maintain the connection with remote machine.

**email:**

type: str, optional

argument path:

simplify\_mdata/model\_devi/  
machine[BohriumContext]/  
remote\_profile/email

Email

**password:**

type: str, optional

argument path:  
 simplify\_mdata/model\_devi/  
 machine[BohriumContext]/  
 remote\_profile/password

Password

**program\_id:**

type: int, alias: *project\_id*  
 argument path:  
 simplify\_mdata/model\_devi/  
 machine[BohriumContext]/  
 remote\_profile/program\_id

Program ID

**retry\_count:**

type: NoneType | int, optional,  
 default: 2  
 argument path:  
 simplify\_mdata/model\_devi/  
 machine[BohriumContext]/  
 remote\_profile/retry\_count

The retry count when a job is terminated

**ignore\_exit\_code:**

type: bool, optional, default: True  
 argument path:  
 simplify\_mdata/model\_devi/  
 machine[BohriumContext]/  
 remote\_profile/  
 ignore\_exit\_code

**The job state will be marked as finished if the exit code is non-zero when set to True. Otherwise,**  
 the job state will be designated as terminated.

**keep\_backup:**

type: bool, optional  
 argument path:  
 simplify\_mdata/model\_devi/  
 machine[BohriumContext]/  
 remote\_profile/keep\_backup

keep download and upload zip

**input\_data:**

type: dict  
 argument path:  
 simplify\_mdata/model\_devi/  
 machine[BohriumContext]/  
 remote\_profile/input\_data

Configuration of job

When `context_type` is set to `LocalContext` (or its aliases `localcontext`, `Local`, `local`):

**remote\_profile:**

type: dict, optional  
argument path: `simplify_mdata/  
model_devi/machine[LocalContext]/  
remote_profile`

The information used to maintain the connection with remote machine. This field is empty for this context.

When `context_type` is set to `LazyLocalContext` (or its aliases `lazylocalcontext`, `LazyLocal`, `lazylocal`):

**remote\_profile:**

type: dict, optional  
argument path: `simplify_mdata/model_devi/  
machine[LazyLocalContext]/  
remote_profile`

The information used to maintain the connection with remote machine. This field is empty for this context.

When `context_type` is set to `SSHContext` (or its aliases `sshcontext`, `SSH`, `ssh`):

**remote\_profile:**

type: dict  
argument path: `simplify_mdata/model_devi/  
machine[SSHContext]/remote_profile`

The information used to maintain the connection with remote machine.

**hostname:**

type: str  
argument path: `simplify_mdata/model_devi/  
machine[SSHContext]/  
remote_profile/hostname`

hostname or ip of ssh connection.

**username:**

type: str  
argument path: `simplify_mdata/model_devi/  
machine[SSHContext]/  
remote_profile/username`

username of target linux system

**password:**

type: str, optional  
 argument path:  
 simplify\_mdata/model\_devi/  
 machine[SSHContext]/  
 remote\_profile/password  
 (deprecated) password of linux system. Please use [SSH keys](#) instead to improve security.

**port:**

type: int, optional, default: 22  
 argument path:  
 simplify\_mdata/model\_devi/  
 machine[SSHContext]/  
 remote\_profile/port  
 ssh connection port.

**key\_filename:**

type: str | NoneType, optional,  
 default: None  
 argument path:  
 simplify\_mdata/model\_devi/  
 machine[SSHContext]/  
 remote\_profile/key\_filename  
 key filename used by ssh connection.  
 If left None, find key in ~/.ssh or use  
 password for login

**passphrase:**

type: str | NoneType, optional,  
 default: None  
 argument path:  
 simplify\_mdata/model\_devi/  
 machine[SSHContext]/  
 remote\_profile/passphrase  
 passphrase of key used by ssh connection

**timeout:**

type: int, optional, default: 10  
 argument path:  
 simplify\_mdata/model\_devi/  
 machine[SSHContext]/  
 remote\_profile/timeout  
 timeout of ssh connection

**totp\_secret:**

type: str | NoneType, optional,  
 default: None

argument path:  
simplify\_mdata/model\_devi/  
machine[SSHContext]/  
remote\_profile/totp\_secret

Time-based one time password secret. It should be a base32-encoded string extracted from the 2D code.

**tar\_compress:**

type: bool, optional, default: True  
argument path:  
simplify\_mdata/model\_devi/  
machine[SSHContext]/  
remote\_profile/tar\_compress

The archive will be compressed in upload and download if it is True. If not, compression will be skipped.

**look\_for\_keys:**

type: bool, optional, default: True  
argument path:  
simplify\_mdata/model\_devi/  
machine[SSHContext]/  
remote\_profile/look\_for\_keys

enable searching for discoverable private key files in ~/.ssh/

When `context_type` is set to `HDFSContext` (or its aliases `hdfscontext`, `HDFS`, `hdfs`):

**remote\_profile:**

type: dict, optional  
argument path:  
simplify\_mdata/model\_devi/  
machine[HDFSContext]/remote\_profile

The information used to maintain the connection with remote machine. This field is empty for this context.

When `context_type` is set to `OpenAPIContext` (or its aliases `openapicontext`, `OpenAPI`, `openapi`):

**remote\_profile:**

type: dict, optional  
argument path: simplify\_mdata/  
model\_devi/machine[OpenAPIContext]/  
remote\_profile

The information used to maintain the connection with remote machine. This field is empty for this context.

**resources:**

type: dict  
 argument path:  
 simplify\_mdata/model\_devi/resources

**number\_node:**

type: int, optional, default: 1  
 argument path: simplify\_mdata/  
 model\_devi/resources/number\_node

The number of node need for each *job*

**cpu\_per\_node:**

type: int, optional, default: 1  
 argument path: simplify\_mdata/  
 model\_devi/resources/cpu\_per\_node

cpu numbers of each node assigned to each job.

**gpu\_per\_node:**

type: int, optional, default: 0  
 argument path: simplify\_mdata/  
 model\_devi/resources/gpu\_per\_node

gpu numbers of each node assigned to each job.

**queue\_name:**

type: str, optional, default: (empty string)  
 argument path: simplify\_mdata/  
 model\_devi/resources/queue\_name

The queue name of batch job scheduler system.

**group\_size:**

type: int  
 argument path: simplify\_mdata/  
 model\_devi/resources/group\_size

The number of *tasks* in a *job*. 0 means infinity.

**custom\_flags:**

type: typing.List[str], optional  
 argument path: simplify\_mdata/  
 model\_devi/resources/custom\_flags

The extra lines pass to job submitting script header

**strategy:**

type: dict, optional  
 argument path: simplify\_mdata/  
 model\_devi/resources/strategy

strategies we use to generation job submitting scripts.

**if\_cuda\_multi\_devices:**

type: bool, optional, default: False  
argument path:  
simplify\_mdata/model\_devi/  
resources/strategy/  
if\_cuda\_multi\_devices

If there are multiple nvidia GPUS on the node, and we want to assign the tasks to different GPUS. If true, dpdispatcher will manually export environment variable CUDA\_VISIBLE\_DEVICES to different task. Usually, this option will be used with Task.task\_need\_resources variable simultaneously.

**ratio\_unfinished:**

type: float, optional, default: 0.0  
argument path: simplify\_mdata/  
model\_devi/resources/  
strategy/ratio\_unfinished

The ratio of *tasks* that can be unfinished.

**customized\_script\_header\_template\_file:**

type: str, optional  
argument path:  
simplify\_mdata/model\_devi/  
resources/strategy/  
customized\_script\_header\_template\_file

The customized template file to generate job submitting script header, which overrides the default file.

**para\_deg:**

type: int, optional, default: 1  
argument path: simplify\_mdata/  
model\_devi/resources/para\_deg

Decide how many tasks will be run in parallel.

**source\_list:**

type: typing.List[str], optional, default: []  
argument path: simplify\_mdata/  
model\_devi/resources/source\_list

The env file to be sourced before the command execution.



**module\_purge:**

type: bool, optional, default: False

argument path: simplify\_mdata/  
model\_devi/resources/module\_purge

Remove all modules on HPC system before  
module load (module\_list)

**module\_unload\_list:**

type: typing.List[str], optional, default:  
[]

argument path:  
simplify\_mdata/model\_devi/  
resources/module\_unload\_list

The modules to be unloaded on HPC system  
before submitting jobs

**module\_list:**

type: typing.List[str], optional, default:  
[]

argument path: simplify\_mdata/  
model\_devi/resources/module\_list

The modules to be loaded on HPC system be-  
fore submitting jobs

**envs:**

type: dict, optional, default: {}

argument path: simplify\_mdata/  
model\_devi/resources/envs

The environment variables to be exported on  
before submitting jobs

**prepend\_script:**

type: typing.List[str], optional, default:  
[]

argument path: simplify\_mdata/  
model\_devi/resources/prepend\_script

Optional script run before jobs submitted.

**append\_script:**

type: typing.List[str], optional, default:  
[]

argument path: simplify\_mdata/  
model\_devi/resources/append\_script

Optional script run after jobs submitted.

**wait\_time:**

type: float | int, optional, default: 0

argument path: `simplify_mdata/  
model_devi/resources/wait_time`

The waiting time in second after a single *task* submitted

Depending on the value of *batch\_type*, different sub args are accepted.

**batch\_type:**

type: str (flag key)

argument path: `simplify_mdata/  
model_devi/resources/batch_type`

possible choices: *Bohrium*, *Fugaku*, *Slurm*,  
*SlurmJobArray*, *Shell*, *LSF*, *PBS*, *SGE*,  
*Torque*, *OpenAPI*, *DistributedShell*

The batch job system type loaded from machine/*batch\_type*.

When *batch\_type* is set to *Bohrium* (or its aliases *bohrium*, *Lebesgue*, *lebesgue*, *DpCloudServer*, *dpcloudserver*):

**kwargs:**

type: dict, optional

argument path:  
`simplify_mdata/model_devi/  
resources[Bohrium]/kwargs`

This field is empty for this batch.

When *batch\_type* is set to *Fugaku* (or its alias *fugaku*):

**kwargs:**

type: dict, optional

argument path: `simplify_mdata/  
model_devi/resources[Fugaku]/kwargs`

This field is empty for this batch.

When *batch\_type* is set to *Slurm* (or its alias *slurm*):

**kwargs:**

type: dict, optional

argument path: `simplify_mdata/  
model_devi/resources[Slurm]/kwargs`

Extra arguments.

**custom\_gpu\_line:**

type: str | NoneType, optional,  
default: None

argument path: `simplify_mdata/  
model_devi/resources[Slurm]/  
kwargs/custom_gpu_line`

Custom GPU configuration, starting with *#SBATCH*

When `batch_type` is set to `SlurmJobArray` (or its alias `slurmjobarray`):

**kwargs:**

type: dict, optional  
 argument path:  
`simplify_mdata/model_devi/  
 resources[SlurmJobArray]/kwargs`

Extra arguments.

**custom\_gpu\_line:**

type: str | NoneType, optional,  
 default: None  
 argument path:  
`simplify_mdata/model_devi/  
 resources[SlurmJobArray]/  
 kwargs/custom_gpu_line`  
 Custom GPU configuration, starting  
 with `#SBATCH`

**slurm\_job\_size:**

type: int, optional, default: 1  
 argument path:  
`simplify_mdata/model_devi/  
 resources[SlurmJobArray]/  
 kwargs/slurm_job_size`  
 Number of tasks in a Slurm job

When `batch_type` is set to `Shell` (or its alias `shell`):

**kwargs:**

type: dict, optional  
 argument path: `simplify_mdata/  
 model_devi/resources[Shell]/kwargs`  
 This field is empty for this batch.

When `batch_type` is set to `LSF` (or its alias `lsf`):

**kwargs:**

type: dict  
 argument path: `simplify_mdata/  
 model_devi/resources[LSF]/kwargs`  
 Extra arguments.

**gpu\_usage:**

type: bool, optional, default: False  
 argument path: `simplify_mdata/  
 model_devi/resources[LSF]/  
 kwargs/gpu_usage`  
 Choosing if GPU is used in the calcu-  
 lation step.

**gpu\_new\_syntax:**

type: bool, optional, default: False  
argument path: simplify\_mdata/  
model\_devi/resources[LSF]/  
kwargs/gpu\_new\_syntax

For LFS >= 10.1.0.3, new option -gpu  
for #BSUB could be used. If False,  
and old syntax would be used.

**gpu\_exclusive:**

type: bool, optional, default: True  
argument path: simplify\_mdata/  
model\_devi/resources[LSF]/  
kwargs/gpu\_exclusive

Only take effect when new syntax en-  
abled. Control whether submit tasks  
in exclusive way for GPU.

**custom\_gpu\_line:**

type: str | NoneType, optional,  
default: None  
argument path: simplify\_mdata/  
model\_devi/resources[LSF]/  
kwargs/custom\_gpu\_line

Custom GPU configuration, starting  
with #BSUB

When `batch_type` is set to PBS (or its alias pbs):

**kwargs:**

type: dict, optional  
argument path: simplify\_mdata/  
model\_devi/resources[PBS]/kwargs

This field is empty for this batch.

When `batch_type` is set to SGE (or its alias sge):

**kwargs:**

type: dict, optional  
argument path: simplify\_mdata/  
model\_devi/resources[SGE]/kwargs

This field is empty for this batch.

When `batch_type` is set to Torque (or its alias torque):

**kwargs:**

type: dict, optional  
argument path: simplify\_mdata/  
model\_devi/resources[Torque]/kwargs

This field is empty for this batch.

When `batch_type` is set to OpenAPI (or its alias openapi):

**kwargs:**

type: dict, optional  
 argument path:  
 simplify\_mdata/model\_devi/  
 resources[OpenAPI]/kwargs

This field is empty for this batch.

When `batch_type` is set to DistributedShell (or its alias distributedshell):

**kwargs:**

type: dict, optional  
 argument path:  
 simplify\_mdata/model\_devi/  
 resources[DistributedShell]/kwargs

This field is empty for this batch.

**user\_forward\_files:**

type: list, optional  
 argument path:  
 simplify\_mdata/model\_devi/user\_forward\_files

Files to be forwarded to the remote machine.

**user\_backward\_files:**

type: list, optional  
 argument path:  
 simplify\_mdata/model\_devi/user\_backward\_files

Files to be backwarded from the remote machine.

**fp:**

type: dict  
 argument path: simplify\_mdata/fp

Parameters of command, machine, and resources for fp

**command:**

type: str  
 argument path: simplify\_mdata/fp/command

Command of a program.

**machine:**

type: dict  
 argument path: simplify\_mdata/fp/machine

**batch\_type:**

type: str  
 argument path: simplify\_mdata/fp/  
 machine/batch\_type

The batch job system type. Option: Fugaku, Slurm, DistributedShell, SGE, PBS, Torque, SlurmJobArray, Shell, Bohrium, LSF, OpenAPI

**local\_root:**

type: str | NoneType  
argument path: simplify\_mdata/fp/  
machine/local\_root

The dir where the tasks and relating files locate. Typically the project dir.

**remote\_root:**

type: str | NoneType, optional  
argument path: simplify\_mdata/fp/  
machine/remote\_root

The dir where the tasks are executed on the remote machine. Only needed when context is not lazy-local.

**clean\_asynchronously:**

type: bool, optional, default: False  
argument path: simplify\_mdata/fp/  
machine/clean\_asynchronously

Clean the remote directory asynchronously after the job finishes.

Depending on the value of *context\_type*, different sub args are accepted.

**context\_type:**

type: str (flag key)  
argument path: simplify\_mdata/fp/  
machine/context\_type  
possible choices: *BohriumContext*,  
*LocalContext*, *LazyLocalContext*,  
*SSHContext*, *HDFSContext*,  
*OpenAPIContext*

The connection used to remote machine. Option: OpenAPIContext, SSHContext, BohriumContext, HDFSContext, LazyLocalContext, LocalContext

When *context\_type* is set to *BohriumContext* (or its aliases *bohriumcontext*, *Bohrium*, *bohrium*, *DpCloudServerContext*, *dpcloudservercontext*, *DpCloudServer*, *dpcloudserver*, *LebesgueContext*, *lebesguecontext*, *Lebesgue*, *lebesgue*):

**remote\_profile:**

type: dict

argument path: simplify\_mdata/fp/  
 machine[BohriumContext]/  
 remote\_profile

The information used to maintain the connection with remote machine.

**email:**

type: str, optional  
 argument path: simplify\_mdata/  
 fp/machine[BohriumContext]/  
 remote\_profile/email

Email

**password:**

type: str, optional  
 argument path: simplify\_mdata/  
 fp/machine[BohriumContext]/  
 remote\_profile/password

Password

**program\_id:**

type: int, alias: *project\_id*  
 argument path: simplify\_mdata/  
 fp/machine[BohriumContext]/  
 remote\_profile/program\_id

Program ID

**retry\_count:**

type: NoneType | int, optional,  
 default: 2  
 argument path: simplify\_mdata/  
 fp/machine[BohriumContext]/  
 remote\_profile/retry\_count

The retry count when a job is terminated

**ignore\_exit\_code:**

type: bool, optional, default: True  
 argument path: simplify\_mdata/  
 fp/machine[BohriumContext]/  
 remote\_profile/  
 ignore\_exit\_code

**The job state will be marked as finished if the exit code is non-zero when set to True. Otherwise, the job state will be designated as terminated.**

**keep\_backup:**

type: bool, optional

```
argument path: simplify_mdata/  
fp/machine[BohriumContext]/  
remote_profile/keep_backup
```

keep download and upload zip

**input\_data:**

```
type: dict  
argument path: simplify_mdata/  
fp/machine[BohriumContext]/  
remote_profile/input_data
```

Configuration of job

When `context_type` is set to `LocalContext` (or its aliases `localcontext`, `Local`, `local`):

**remote\_profile:**

```
type: dict, optional  
argument path: simplify_mdata/fp/  
machine[LocalContext]/  
remote_profile
```

The information used to maintain the connection with remote machine. This field is empty for this context.

When `context_type` is set to `LazyLocalContext` (or its aliases `lazylocalcontext`, `LazyLocal`, `lazylocal`):

**remote\_profile:**

```
type: dict, optional  
argument path: simplify_mdata/fp/  
machine[LazyLocalContext]/  
remote_profile
```

The information used to maintain the connection with remote machine. This field is empty for this context.

When `context_type` is set to `SSHContext` (or its aliases `sshcontext`, `SSH`, `ssh`):

**remote\_profile:**

```
type: dict  
argument path: simplify_mdata/fp/  
machine[SSHContext]/remote_profile
```

The information used to maintain the connection with remote machine.

**hostname:**

```
type: str  
argument path: simplify_mdata/  
fp/machine[SSHContext]/  
remote_profile/hostname
```

hostname or ip of ssh connection.



**username:**

type: str  
argument path: simplify\_mdata/  
fp/machine[SSHContext]/  
remote\_profile/username  
username of target linux system

**password:**

type: str, optional  
argument path: simplify\_mdata/  
fp/machine[SSHContext]/  
remote\_profile/password  
(deprecated) password of linux system. Please use [SSH keys](#) instead to improve security.

**port:**

type: int, optional, default: 22  
argument path: simplify\_mdata/  
fp/machine[SSHContext]/  
remote\_profile/port  
ssh connection port.

**key\_filename:**

type: str | NoneType, optional,  
default: None  
argument path: simplify\_mdata/  
fp/machine[SSHContext]/  
remote\_profile/key\_filename  
key filename used by ssh connection.  
If left None, find key in ~/.ssh or use  
password for login

**passphrase:**

type: str | NoneType, optional,  
default: None  
argument path: simplify\_mdata/  
fp/machine[SSHContext]/  
remote\_profile/passphrase  
passphrase of key used by ssh connection

**timeout:**

type: int, optional, default: 10  
argument path: simplify\_mdata/  
fp/machine[SSHContext]/  
remote\_profile/timeout  
timeout of ssh connection

**totp\_secret:**

type: str | NoneType, optional,  
default: None

argument path: `simplify_mdata/  
fp/machine[SSHContext]/  
remote_profile/totp_secret`

Time-based one time password secret. It should be a base32-encoded string extracted from the 2D code.

**tar\_compress:**

type: bool, optional, default: True  
argument path: `simplify_mdata/  
fp/machine[SSHContext]/  
remote_profile/tar_compress`

The archive will be compressed in upload and download if it is True. If not, compression will be skipped.

**look\_for\_keys:**

type: bool, optional, default: True  
argument path: `simplify_mdata/  
fp/machine[SSHContext]/  
remote_profile/look_for_keys`

enable searching for discoverable private key files in `~/.ssh/`

When `context_type` is set to `HDFSContext` (or its aliases `hdfscontext`, `HDFS`, `hdfs`):

**remote\_profile:**

type: dict, optional  
argument path: `simplify_mdata/fp/  
machine[HDFSContext]/remote_profile`

The information used to maintain the connection with remote machine. This field is empty for this context.

When `context_type` is set to `OpenAPIContext` (or its aliases `openapicontext`, `OpenAPI`, `openapi`):

**remote\_profile:**

type: dict, optional  
argument path: `simplify_mdata/fp/  
machine[OpenAPIContext]/  
remote_profile`

The information used to maintain the connection with remote machine. This field is empty for this context.

**resources:**

type: dict  
argument path: `simplify_mdata/fp/resources`

**number\_node:**

type: int, optional, default: 1  
 argument path: simplify\_mdata/fp/  
 resources/number\_node

The number of node need for each *job*

**cpu\_per\_node:**

type: int, optional, default: 1  
 argument path: simplify\_mdata/fp/  
 resources/cpu\_per\_node

cpu numbers of each node assigned to each job.

**gpu\_per\_node:**

type: int, optional, default: 0  
 argument path: simplify\_mdata/fp/  
 resources/gpu\_per\_node

gpu numbers of each node assigned to each job.

**queue\_name:**

type: str, optional, default: (empty string)  
 argument path: simplify\_mdata/fp/  
 resources/queue\_name

The queue name of batch job scheduler system.

**group\_size:**

type: int  
 argument path: simplify\_mdata/fp/  
 resources/group\_size

The number of *tasks* in a *job*. 0 means infinity.

**custom\_flags:**

type: typing.List[str], optional  
 argument path: simplify\_mdata/fp/  
 resources/custom\_flags

The extra lines pass to job submitting script header

**strategy:**

type: dict, optional  
 argument path: simplify\_mdata/fp/  
 resources/strategy

strategies we use to generation job submitting scripts.

**if\_cuda\_multi\_devices:**

type: bool, optional, default: False

argument path: `simplify_mdata/  
fp/resources/strategy/  
if_cuda_multi_devices`

If there are multiple nvidia GPUS on the node, and we want to assign the tasks to different GPUS. If true, `dpdispatcher` will manually export environment variable `CUDA_VISIBLE_DEVICES` to different task. Usually, this option will be used with `Task.task_need_resources` variable simultaneously.

**ratio\_unfinished:**

type: float, optional, default: 0.0

argument path:

`simplify_mdata/fp/resources/  
strategy/ratio_unfinished`

The ratio of *tasks* that can be unfinished.

**customized\_script\_header\_template\_file:**

type: str, optional

argument path: `simplify_mdata/  
fp/resources/strategy/  
customized_script_header_template_file`

The customized template file to generate job submitting script header, which overrides the default file.

**para\_deg:**

type: int, optional, default: 1

argument path: `simplify_mdata/fp/  
resources/para_deg`

Decide how many tasks will be run in parallel.

**source\_list:**

type: `typing.List[str]`, optional, default: []

argument path: `simplify_mdata/fp/  
resources/source_list`

The env file to be sourced before the command execution.

**module\_purge:**

type: bool, optional, default: False

argument path: `simplify_mdata/fp/  
resources/module_purge`

Remove all modules on HPC system before module load (`module_list`)

**module\_unload\_list:**

type: `typing.List[str]`, optional, default: `[]`

argument path: `simplify_mdata/fp/resources/module_unload_list`

The modules to be unloaded on HPC system before submitting jobs

**module\_list:**

type: `typing.List[str]`, optional, default: `[]`

argument path: `simplify_mdata/fp/resources/module_list`

The modules to be loaded on HPC system before submitting jobs

**envs:**

type: `dict`, optional, default: `{}`

argument path: `simplify_mdata/fp/resources/envs`

The environment variables to be exported on before submitting jobs

**prepend\_script:**

type: `typing.List[str]`, optional, default: `[]`

argument path: `simplify_mdata/fp/resources/prepend_script`

Optional script run before jobs submitted.

**append\_script:**

type: `typing.List[str]`, optional, default: `[]`

argument path: `simplify_mdata/fp/resources/append_script`

Optional script run after jobs submitted.

**wait\_time:**

type: `float | int`, optional, default: `0`

argument path: `simplify_mdata/fp/resources/wait_time`

The waiting time in second after a single *task* submitted

Depending on the value of *batch\_type*, different sub args are accepted.

**batch\_type:**

type: str (flag key)  
argument path: `simplify_mdata/fp/resources/batch_type`  
possible choices: *Bohrium*, *Fugaku*, *Slurm*,  
*SlurmJobArray*, *Shell*, *LSF*, *PBS*, *SGE*,  
*Torque*, *OpenAPI*, *DistributedShell*  
The batch job system type loaded from machine/`batch_type`.

When `batch_type` is set to *Bohrium* (or its aliases *bohrium*, *Lebesgue*, *lebesgue*, *DpCloudServer*, *dpcloudserver*):

**kwargs:**

type: dict, optional  
argument path: `simplify_mdata/fp/resources[Bohrium]/kwargs`  
This field is empty for this batch.

When `batch_type` is set to *Fugaku* (or its alias *fugaku*):

**kwargs:**

type: dict, optional  
argument path: `simplify_mdata/fp/resources[Fugaku]/kwargs`  
This field is empty for this batch.

When `batch_type` is set to *Slurm* (or its alias *slurm*):

**kwargs:**

type: dict, optional  
argument path: `simplify_mdata/fp/resources[Slurm]/kwargs`  
Extra arguments.

**custom\_gpu\_line:**

type: str | NoneType, optional,  
default: None  
argument path: `simplify_mdata/fp/resources[Slurm]/kwargs/custom_gpu_line`  
Custom GPU configuration, starting with `#SBATCH`

When `batch_type` is set to *SlurmJobArray* (or its alias *slurmjobarray*):

**kwargs:**

type: dict, optional

argument path: `simplify_mdata/fp/resources[SlurmJobArray]/kwargs`

Extra arguments.

**custom\_gpu\_line:**

type: `str | NoneType`, optional,  
default: `None`

argument path: `simplify_mdata/fp/resources[SlurmJobArray]/kwargs/custom_gpu_line`

Custom GPU configuration, starting with `#SBATCH`

**slurm\_job\_size:**

type: `int`, optional, default: `1`

argument path: `simplify_mdata/fp/resources[SlurmJobArray]/kwargs/slurm_job_size`

Number of tasks in a Slurm job

When `batch_type` is set to `Shell` (or its alias `shell`):

**kwargs:**

type: `dict`, optional

argument path: `simplify_mdata/fp/resources[Shell]/kwargs`

This field is empty for this batch.

When `batch_type` is set to `LSF` (or its alias `lsf`):

**kwargs:**

type: `dict`

argument path: `simplify_mdata/fp/resources[LSF]/kwargs`

Extra arguments.

**gpu\_usage:**

type: `bool`, optional, default: `False`

argument path: `simplify_mdata/fp/resources[LSF]/kwargs/gpu_usage`

Choosing if GPU is used in the calculation step.

**gpu\_new\_syntax:**

type: `bool`, optional, default: `False`

argument path: `simplify_mdata/fp/resources[LSF]/kwargs/gpu_new_syntax`

For LFS  $\geq$  10.1.0.3, new option `-gpu` for `#BSUB` could be used. If False, and old syntax would be used.

**gpu\_exclusive:**

type: bool, optional, default: True  
argument path: `simplify_mdata/  
fp/resources[LSF]/kwargs/  
gpu_exclusive`

Only take effect when new syntax enabled. Control whether submit tasks in exclusive way for GPU.

**custom\_gpu\_line:**

type: str | NoneType, optional,  
default: None  
argument path: `simplify_mdata/  
fp/resources[LSF]/kwargs/  
custom_gpu_line`

Custom GPU configuration, starting with `#BSUB`

When `batch_type` is set to PBS (or its alias `pbs`):

**kwargs:**

type: dict, optional  
argument path: `simplify_mdata/fp/  
resources[PBS]/kwargs`

This field is empty for this batch.

When `batch_type` is set to SGE (or its alias `sge`):

**kwargs:**

type: dict, optional  
argument path: `simplify_mdata/fp/  
resources[SGE]/kwargs`

This field is empty for this batch.

When `batch_type` is set to Torque (or its alias `torque`):

**kwargs:**

type: dict, optional  
argument path: `simplify_mdata/fp/  
resources[Torque]/kwargs`

This field is empty for this batch.

When `batch_type` is set to OpenAPI (or its alias `openapi`):

**kwargs:**

type: dict, optional



argument path: `simplify_mdata/fp/`  
resources[OpenAPI]/kwargs

This field is empty for this batch.

When `batch_type` is set to `DistributedShell` (or its alias `distributedshell`):

**kwargs:**

type: dict, optional  
argument path: `simplify_mdata/fp/`  
resources[DistributedShell]/kwargs

This field is empty for this batch.

**user\_forward\_files:**

type: list, optional  
argument path:  
`simplify_mdata/fp/user_forward_files`

Files to be forwarded to the remote machine.

**user\_backward\_files:**

type: list, optional  
argument path:  
`simplify_mdata/fp/user_backward_files`

Files to be backwarded from the remote machine.



## **AUTO TEST**

### **7.1 Autotest Overview: Autotest for Deep Generator**

Suppose that we have a potential (can be DFT, DP, MEAM...), `autotest` helps us automatically calculate  $M$  properties on  $N$  configurations. The folder where the `autotest` runs is called the working directory of `autotest`. Different potentials should be tested in different working directories.

A property is tested in three steps: `make`, `run` and `post`. `make` prepares all computational tasks that are needed to calculate the property. For example to calculate EOS, `make` prepares a series of tasks, each of which has a scaled configuration with certain volume, and all necessary input files necessary for starting a VASP, ABACUS, or LAMMPS calculations. `run` sends all the computational tasks to remote computational resources defined in a machine configuration file like `machine.json`, and automatically collects the results when remote calculations finish. `post` calculates the desired property from the collected results.

#### **7.1.1 Relaxation**

The relaxation of a structure should be carried out before calculating all other properties:

```
dpngen autotest make relax.json
dpngen autotest run relax.json machine.json
dpngen autotest post relax.json
```

If, for some reasons, the main program terminated at stage `run`, one can easily restart with the same command. `relax.json` is the parameter file. An example for `deepmd` relaxation is given as:

```
{
  "structures": ["confs/mp-*"],
  "interaction": {
    "type": "deepmd",
    "model": "frozen_model.pb",
    "type_map": {"Al": 0, "Mg": 1}
  },
  "relaxation": {}
}
```

where the key `structures` provides the structures to relax. `interaction` is provided with `deepmd`, and other options are `vasp`, `abacus`, `meam`...

## 7.1.2 Task type

There are now six task types implemented in the package: `vasp`, `abacus`, `deepmd`, `meam`, `eam_fs`, and `eam_alloy`. An `inter.json` file in json format containing the interaction parameters will be written in the directory of each task after make. We give input examples of the `interaction` part for each type below:

### VASP:

The default of `potcar_prefix` is `""`.

```
"interaction": {
  "type": "vasp",
  "incar": "vasp_input/INCAR",
  "potcar_prefix": "vasp_input",
  "potcars": {"Al": "POTCAR.al", "Mg": "POTCAR.mg"}
}
```

### ABACUS:

The default of `potcar_prefix` is `""`. The path of `potcars/orb_files/deepks_desc` is `potcar_prefix + potcars/orb_files/deepks_desc/deepks_model`.

```
"interaction": {
  "type": "abacus",
  "incar": "abacus_input/INPUT",
  "potcar_prefix": "abacus_input",
  "potcars": {"Al": "pseudo_potential.al", "Mg": "pseudo_potential.
↪mg"},
  "orb_files": {"Al": "numerical_orb.al", "Mg": "numerical_orb.mg"},
  "atom_masses": {"Al": 26.9815, "Mg": 24.305},
  "deepks_desc": "jle.orb",
  "deepks_model": "model.ptg"
}
```

### deepmd:

Only 1 model can be used in autotest in one working directory.

```
"interaction": {
  "type": "deepmd",
  "model": "frozen_model.pb",
  "type_map": {"Al": 0, "Mg": 1}
}
```

### meam:

Please make sure the [USER-MEAMC](#) package has already been installed in LAMMPS.

```
"interaction": {
  "type": "meam",
  "model": ["meam.lib", "AlMg.meam"],
  "type_map": {"Al": 1, "Mg": 2}
}
```

### eam\_fs & eam\_alloy:

Please make sure the [MANYBODY](#) package has already been installed in LAMMPS

```

"interaction": {
  "type":          "eam_fs (eam_alloy)",
  "model":         "AlMg.eam.fs (AlMg.eam.alloy)",
  "type_map":      {"Al": 1, "Mg": 2}
}

```

### 7.1.3 Property type

Now the supported property types are `eos`, `elastic`, `vacancy`, `interstitial`, `surface`, and `gamma`. Before property tests, `relaxation` should be done first or the relaxation results should be present in the corresponding directory `confs/mp-*/relaxation/relax_task`. A file named `task.json` in json format containing the property parameter will be written in the directory of each task after make step. Multiple property tests can be performed simultaneously.

## 7.2 Make run and post

There are three operations in auto test package, namely `make`, `run`, and `post`. Here we take `eos` property as an example for property type.

### 7.2.1 Make

The INCAR, POSCAR, POTCAR input files for VASP or `in.lammps`, `conf.lmp`, and the interatomic potential files for LAMMPS will be generated in the directory `confs/mp-*/relaxation/relax_task` for relaxation or `confs/mp-*/eos_00/task.[0-9]*[0-9]` for EOS. The `machine.json` file is not needed for make. Example:

```
dpngen autotest make relaxation.json
```

### 7.2.2 Run

The jobs would be dispatched according to the parameter in `machine.json` file and the calculation results would be sent back. Example:

```
dpngen autotest run relaxation.json machine.json
```

### 7.2.3 Post

The post process of calculation results would be performed. `result.json` in json format will be generated in `confs/mp-*/relaxation/relax_task` for relaxation and `result.json` in json format and `result.out` in txt format in `confs/mp-*/eos_00` for EOS. The `machine.json` file is also not needed for post. Example:

```
dpngen autotest post relaxation.json
```

## 7.3 Relaxation

### 7.3.1 Relaxation get started and input examples

The relaxation of a structure should be carried out before calculating all other properties.

First, we need input parameter file and we name it `relax.json` here. All the relaxation calculations should be taken either by VASP, ABACUS, or LAMMPS. Here are two input examples for VASP and LAMMPS respectively.

An example of the input file for relaxation by VASP:

```
{
  "structures":          ["confs/std-*"],
  "interaction": {
    "type":              "vasp",
    "incar":              "vasp_input/INCAR",
    "potcar_prefix":      "vasp_input",
    "potcars":            {"Al": "POTCAR.al"}
  },
  "relaxation": {
    "cal_type":           "relaxation",
    "cal_setting": {
      "relax_pos":        true,
      "relax_shape":      true,
      "relax_vol":        true,
      "ediff":            1e-6,
      "ediffg":           -0.01,
      "encut":            650,
      "kspacing":         0.1,
      "kgamma":           false}
    }
}
```

Key words	data ture	struc- ture	example		description
<b>structures</b>	List of String		["confs/std-*"]		path of different structures
<b>interaction</b>	Dict		See above		description of the task type and atomic interaction
<b>type</b>	String		"vasp"		task type
<b>incar</b>	String		"vasp_input/INCAR"		path for INCAR file in vasp
pot- car_prefix	String		"vasp_input"		prefix of path for POTCAR file in vasp, default = ""
<b>potcars</b>	Dict		{ "Al": CAR.al" }	"POT-	key is element type and value is potcar name
<b>relaxation</b>	Dict		See above		calculation type and setting for relaxation
cal_type	String		"relaxation" "static"	or	calculation type
cal_setting	Dict		See above		calculation setting
relax_pos	Boolean		true		relax atomic position or not, default = true for relaxation
relax_shape	Boolean		true		relax box shape or not, default = true for relaxation
relax_vol	Boolean		true		relax box volume or not, default = true for relaxation
ediff	Float		1e-6		set EDIFF parameter in INCAR files
ediffg	Float		-0.01		set EDIFFG parameter in INCAR files
encut	Int		650		set encut parameter in INCAR files
kspacing	Float		0.1		set KSPACING parameter in INCAR files
kgamma	Boolean		false		set KGAMMA parameter in INCAR files

An example of the input file for relaxation by LAMMPS:

```
{
  "structures":      ["confs/std-*"],
  "interaction": {
    "type":          "deepmd",
    "model":          "frozen_model.pb",
    "in_lammps":      "lammps_input/in.lammps",
    "type_map":       {"Al": 0}
  },
  "relaxation": {
    "cal_setting": {"etol": 0,
                  "ftol": 1e-10,
                  "maxiter": 5000,
                  "maximal": 500000}
  }
}
```

Other key words different from vasp:

Key words	data structure	example	description
<b>model</b>	String or List of String	“frozen_model.pb”	model file for atomic interaction
<b>in_lammps</b>	String	“lammps_input/in.lan	input file for lammps commands
<b>type_map</b>	Dict	{“Al”: 0}	key is element type and value is type number. DP starts from 0, others starts from 1
etol	Float	0	stopping tolerance for energy
ftol	Float	1e-10	stopping tolerance for force
maxiter	Int	5000	max iterations of minimizer
maxeval	Int	500000	max number of force/energy evaluations

For LAMMPS relaxation and all the property calculations, **package will help to generate in.lammps file for user automatically** according to the property type. We can also make the final changes in the `minimize` setting (`minimize etol ftol maxiter maxeval`) in `in.lammps`. In addition, users can apply the input file for lammps commands in the interaction part. For further information of the LAMMPS relaxation, we refer users to [minimize command](#).

### 7.3.2 Relaxation make

The list of the directories storing structures are `["confs/std-*"]` in the previous example. For single element system, if POSCAR doesn't exist in the directories: `std-fcc`, `std-hcp`, `std-dhcp`, `std-bcc`, `std-diamond`, and `std-sc`, the package will automatically generate the standard crystal structures **fcc**, **hcp**, **dhcp**, **bcc**, **diamond**, and **sc** in the corresponding directories, respectively. In other conditions and for multi-component system (more than 1), if POSCAR doesn't exist, the package will terminate and print the error **“no configuration for autotest”**.

#### VASP relaxation

Take the input example of Al in the previous section, when we do `make` as follows:

```
dpngen autotest make relaxation.json
```

the following files would be generated:

```
tree confs/std-fcc/relaxation/
```

```
confs/std-fcc/relaxation/
|-- INCAR
|-- POTCAR
|-- relax_task
    |-- INCAR -> ../INCAR
    |-- inter.json
    |-- KPOINTS
    |-- POSCAR -> ../../POSCAR
    |-- POTCAR -> ../POTCAR
    |-- task.json
```

`inter.json` records the information in the interaction dictionary and `task.json` records the information in the relaxation dictionary.



## LAMMPS relaxation

```
dpngen autotest make relaxation.json
tree confs/std-fcc/
```

the output would be:

```
confs/std-fcc/
|-- POSCAR
|-- relaxation
|   |-- frozen_model.pb -> ../../frozen_model.pb
|   |-- in.lammps
|   |-- relax_task
|       |-- conf.lmp
|       |-- frozen_model.pb -> ../frozen_model.pb
|       |-- in.lammps -> ../in.lammps
|       |-- inter.json
|       |-- POSCAR -> ../../POSCAR
|       |-- task.json
```

the `conf.lmp` is the input configuration and `in.lammps` is the input command file for lammps.

**in.lammps:** the package would generate the file `confs/mp-*/relaxation/in.lammps` as follows and we refer the user to the further information of `fix box/relax` function in lammps:

```
clear
units                metal
dimension            3
boundary             p p p
atom_style           atomic
box                  tilt large
read_data            conf.lmp
mass                 1 26.982
neigh_modify         every 1 delay 0 check no
pair_style deepmd    frozen_model.pb
pair_coeff
compute              mype all pe
thermo               100
thermo_style         custom step pe pxx pyy pzz pxy pxz pyz lx ly lz vol c_mype
dump                 1 all custom 100 dump.relax id type xs ys zs fx fy fz
min_style            cg
fix                  1 all box/relax iso 0.0
minimize             0 1.000000e-10 5000 500000
fix                  1 all box/relax aniso 0.0
minimize             0 1.000000e-10 5000 500000
variable             N equal count(all)
variable             V equal vol
variable             E equal "c_mype"
variable             tmlpx equal lx
variable             tmlpy equal ly
variable             Pxx equal pxx
variable             Pyy equal pyy
variable             Pzz equal pzz
variable             Pxy equal pxy
```

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```

variable      Pxz equal pxz
variable      Pyz equal pyz
variable      Epa equal ${E}/${N}
variable      Vpa equal ${V}/${N}
variable      AA equal (${tmplx}*${tmply})
print "All done"
print "Total number of atoms = ${N}"
print "Final energy per atoms = ${Epa}"
print "Final volume per atoms = ${Vpa}"
print "Final Base area = ${AA}"
print "Final Stress (xx yy zz xy xz yz) = ${Pxx} ${Pyy} ${Pzz} ${Pxy} ${Pxz} ${Pyz}"

```

If user provides lammps input command file `in.lammps`, the `thermo_style` and `dump` commands should be the same as the above file.

**interatomic potential model:** the `frozen_model.pb` in `confs/mp-*/relaxation` would link to the `frozen_model.pb` file given in the input.

### 7.3.3 Relaxation run

The work path of each task should be in the form like `confs/mp-*/relaxation` and all task is in the form like `confs/mp-*/relaxation/relax_task`.

The `machine.json` file should be applied in this process and the machine parameters (eg. GPU or CPU) are determined according to the task type (VASP or LAMMPS). Then in each work path, the corresponding tasks would be submitted and the results would be sent back through [make\\_dispatcher](#).

Take `deepmd` run for example:

```

nohup dpngen autotest run relaxation.json machine-ali.json > run.result 2>&1 &
tree confs/std-fcc/relaxation/

```

the output would be:

```

confs/std-fcc/relaxation/
|-- frozen_model.pb -> ../../../../frozen_model.pb
|-- in.lammps
|-- jr.json
|-- relax_task
|   |-- conf.lmp
|   |-- dump.relax
|   |-- frozen_model.pb -> ../frozen_model.pb
|   |-- in.lammps -> ../in.lammps
|   |-- inter.json
|   |-- log.lammps
|   |-- outlog
|   |-- POSCAR -> ../../POSCAR
|-- task.json

```

`dump.relax` is the file storing configurations and `log.lammps` is the output file for lammps.

### 7.3.4 Relaxation post

Take deepmd post for example:

```
dpngen autotest post relaxation.json
tree confs/std-fcc/relaxation/
```

the output will be:

```
confs/std-fcc/relaxation/
|-- frozen_model.pb -> ../../../../frozen_model.pb
|-- in.lammps
|-- jr.json
|-- relax_task
|   |-- conf.lmp
|   |-- CONTCAR
|   |-- dump.relax
|   |-- frozen_model.pb -> ../frozen_model.pb
|   |-- in.lammps -> ../in.lammps
|   |-- inter.json
|   |-- log.lammps
|   |-- outlog
|   |-- POSCAR -> ../../POSCAR
|   |-- result.json
|   |-- task.json
```

result.json stores the box cell, coordinates, energy, force, virial,... information of each frame in the relaxation trajectory and CONTCAR is the final equilibrium configuration.

result.json:

```
{
  "@module": "dpdata.system",
  "@class": "LabeledSystem",
  "data": {
    "atom_nums": [
      1
    ],
    "atom_names": [
      "Al"
    ],
    "atom_types": {
      "@module": "numpy",
      "@class": "array",
      "dtype": "int64",
      "data": [
        0
      ]
    },
    "orig": {
      "@module": "numpy",
      "@class": "array",
      "dtype": "int64",
      "data": [
```

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```

        0,
        0,
        0
    ]
},
"cells": {
    "@module": "numpy",
    "@class": "array",
    "dtype": "float64",
    "data": [
        [
            [
                2.8637824638,
                0.0,
                0.0
            ],
            [
                1.4318912319,
                2.4801083646,
                0.0
            ],
            [
                1.4318912319,
                0.8267027882,
                2.3382685902
            ]
        ],
        [
            [
                2.8549207998018438,
                0.0,
                0.0
            ],
            [
                1.4274603999009239,
                2.472433938457684,
                0.0
            ],
            [
                1.4274603999009212,
                0.8241446461525599,
                2.331033071844216
            ]
        ],
        [
            [
                2.854920788303194,
                0.0,
                0.0
            ],
            [
                1.427460394144466,

```

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```

                2.472433928487206,
                0.0
            ],
            [
                1.427460394154763,
                0.8241446428350139,
                2.331033062460779
            ]
        ]
    ],
    "coords": {
        "@module": "numpy",
        "@class": "array",
        "dtype": "float64",
        "data": [
            [
                [
                    0.0,
                    0.0,
                    0.0
                ]
            ],
            [
                [
                    5.709841595683707e-25,
                    -4.3367974740910857e-19,
                    0.0
                ]
            ],
            [
                [
                    -8.673606219968035e-19,
                    8.673619637565944e-19,
                    8.673610853102186e-19
                ]
            ]
        ]
    },
    "energies": {
        "@module": "numpy",
        "@class": "array",
        "dtype": "float64",
        "data": [
            -3.745029,
            -3.7453815,
            -3.7453815
        ]
    },
    "forces": {
        "@module": "numpy",
        "@class": "array",

```

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```

    "dtype": "float64",
    "data": [
        [
            [
                0.0,
                -6.93889e-18,
                -3.46945e-18
            ]
        ],
        [
            [
                1.38778e-17,
                6.93889e-18,
                -1.73472e-17
            ]
        ],
        [
            [
                1.38778e-17,
                1.73472e-17,
                -4.51028e-17
            ]
        ]
    ],
    "virials": {
        "@module": "numpy",
        "@class": "array",
        "dtype": "float64",
        "data": [
            [
                [
                    -0.07534992071654338,
                    1.2156615579052586e-17,
                    1.3904892126132796e-17
                ],
                [
                    1.2156615579052586e-17,
                    -0.07534992071654338,
                    4.61571024026576e-12
                ],
                [
                    1.3904892126132796e-17,
                    4.61571024026576e-12,
                    -0.07534992071654338
                ]
            ],
            [
                [
                    -9.978994290457664e-08,
                    -3.396452753975288e-15,
                    8.785831629151552e-16
                ]
            ]
        ]
    }

```

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```

    ],
    [
        -3.396452753975288e-15,
        -9.991375413666671e-08,
        5.4790751628409565e-12
    ],
    [
        8.785831629151552e-16,
        5.4790751628409565e-12,
        -9.973497959053003e-08
    ]
],
[
    [
        1.506940521266962e-11,
        1.1152016233536118e-11,
        -8.231900529157644e-12
    ],
    [
        1.1152016233536118e-11,
        -6.517665029355618e-11,
        -6.33706710415926e-12
    ],
    [
        -8.231900529157644e-12,
        -6.33706710415926e-12,
        5.0011471096530724e-11
    ]
]
]
},
"stress": {
    "@module": "numpy",
    "@class": "array",
    "dtype": "float64",
    "data": [
        [
            -7.2692250000000005,
            1.1727839e-15,
            1.3414452e-15
        ],
        [
            1.1727839e-15,
            -7.2692250000000005,
            4.4529093000000003e-10
        ],
        [
            1.3414452e-15,
            4.4529093000000003e-10,
            -7.2692250000000005
        ]
    ]
}

```

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```
[
    [
        -9.71695e-06,
        -3.3072633e-13,
        8.5551193e-14
    ],
    [
        -3.3072633e-13,
        -9.7290060000000001e-06,
        5.3351969e-10
    ],
    [
        8.5551193e-14,
        5.3351969e-10,
        -9.711598e-06
    ]
],
[
    [
        1.4673689e-09,
        1.0859169e-09,
        -8.0157343e-10
    ],
    [
        1.0859169e-09,
        -6.3465139e-09,
        -6.1706584e-10
    ],
    [
        -8.0157343e-10,
        -6.1706584e-10,
        4.8698191e-09
    ]
]
}
```

## 7.4 Property

### 7.4.1 Property get started and input examples

Here we take deepmd for example and the input file for other task types is similar.

```
{
  "structures":      ["confs/std-*"],
  "interaction": {
    "type":          "deepmd",
    (continues on next page)
```

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```

    "model":      "frozen_model.pb",
    "type_map":   {"Al": 0}
  },
  "properties": [
    {
      "type":      "eos",
      "vol_start": 0.9,
      "vol_end":   1.1,
      "vol_step":  0.01
    },
    {
      "type":      "elastic",
      "norm_deform": 1e-2,
      "shear_deform": 1e-2
    },
    {
      "type":      "vacancy",
      "supercell":  [3, 3, 3],
      "start_confs_path": "../vasp/confs"
    },
    {
      "type":      "interstitial",
      "supercell":  [3, 3, 3],
      "insert_ele": ["Al"],
      "conf_filters": {"min_dist": 1.5},
      "cal_setting": {"input_prop": "lammps_input/lammps_high"}
    },
    {
      "type":      "surface",
      "min_slab_size": 10,
      "min_vacuum_size": 11,
      "max_miller":   2,
      "cal_type":     "static"
    },
    {
      "type": "gamma",
      "lattice_type": "fcc",
      "miller_index": [1, 1, 1],
      "displace_direction": [1, 1, 0],
      "supercell_size": [1, 1, 10],
      "min_vacuum_size": 10,
      "add_fix": ["true", "true", "false"],
      "n_steps": 20
    }
  ]
}

```

Universal key words for properties

Key words	data structure	example	description
<b>type</b>	String	“eos”	property type
<b>skip</b>	Boolean	true	whether to skip current property or not
<b>start_conf_path</b>	String	“../vasp/confs”	start from the equilibrium configuration in other path only for the current property type
<b>cal_setting[“input_prop”]</b>	String	“lammps_input/lamr	input commands file
<b>cal_setting[“overwrite_i”]</b>	Dict		overwrite the interaction in the <code>interaction</code> part only for the current property type

other parameters in `cal_setting` and `cal_type` in `relaxation` also apply in `property`.

Key words for **EOS**

Key words	data structure	example	description
<b>vol_start</b>	Float	0.9	the starting volume related to the equilibrium structure
<b>vol_end</b>	Float	1.1	the biggest volume related to the equilibrium structure
<b>vol_step</b>	Float	0.01	the volume increment related to the equilibrium structure
<b>vol_abs</b>	Boolean	false	whether to treat <code>vol_start</code> , <code>vol_end</code> and <code>vol_step</code> as absolute volume or not (as relative volume), default = false

Key words for **Elastic**

Key words	data structure	example	description
<b>norm_deform</b>	Float	1e-2	deformation in xx, yy, zz, default = 1e-2
<b>shear_deform</b>	Float	1e-2	deformation in other directions, default = 1e-2

Key words for **Vacancy**

Key words	data structure	example	description
<b>supercell</b>	List of Int	[3,3,3]	the supercell to be constructed, default = [1,1,1]

Key words for **Interstitial**

Key words	data structure	example	description
<b>in-sert_ele</b>	List of String	[“Al”]	the element to be inserted
<b>supercell</b>	List of Int	[3,3,3]	the supercell to be constructed, default = [1,1,1]
<b>conf_filters</b>	Dict	“min_dist”: 1.5	filter out the undesirable configuration
<b>bcc_self</b>	Boolean	false	whether to do the self-interstitial calculations for bcc structures, default = false

Key words for **Surface**

Key words	data structure	example	description
<b>min_slab_size</b>	Int	10	minimum size of slab thickness
<b>min_vacuum_size</b>	Int	11	minimum size of vacuum width
<b>pert_xz</b>	Float	0.01	perturbation through xz direction used to compute surface energy, default = 0.01
<b>max_miller</b>	Int	2	the maximum miller index, default = 2

Key words for **Gamma**

Key words	data structure	example	description
<b>lattice_type</b>	String	“fcc”	“bcc” or “fcc” at this stage
<b>miller_index</b>	List of Int	[1,1,1]	slip plane for gamma-line calculation
<b>displace_direction</b>	List of Int	[1,1,0]	slip direction for gamma-line calculation
<b>supercell_size</b>	List of Int	[1,1,10]	the supercell to be constructed, default = [1,1,5]
<b>min_vacuum_size</b>	Int or Float	10	minimum size of vacuum width, default = 20
<b>add_fix</b>	List of String	['true','true','false']	whether to fix atoms in the direction, default = ['true','true','false'] (standard method)
<b>n_steps</b>	Int	20	Number of points for gamma-line calculation, default = 10

## 7.4.2 Property make

```
dpngen autotest make property.json
```

## EOS output:

```
confs/std-fcc/eos_00/
|-- frozen_model.pb -> ../../../../frozen_model.pb
|-- task.0000000
|   |-- conf.lmp
|   |-- eos.json
|   |-- frozen_model.pb -> ../frozen_model.pb
|   |-- in.lammps
|   |-- inter.json
|   |-- POSCAR
|   |-- POSCAR.orig -> ../../relaxation/relax_task/CONTCAR
|   |-- task.json
|-- task.0000001
|   |-- conf.lmp
|   |-- eos.json
|   |-- frozen_model.pb -> ../frozen_model.pb
|   |-- in.lammps
|   |-- inter.json
|   |-- POSCAR
|   |-- POSCAR.orig -> ../../relaxation/relax_task/CONTCAR
```

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```
| -- task.json
...
|-- task.000019
| -- conf.lmp
| -- eos.json
| -- frozen_model.pb -> ../frozen_model.pb
| -- in.lammps
| -- inter.json
| -- POSCAR
| -- POSCAR.orig -> ../../relaxation/relax_task/CONTCAR
|-- task.json
```

eos.json records the volume and scale of the corresponding task.

#### Elastic output:

```
confs/std-fcc/elastic_00/
|-- equi.stress.json
|-- frozen_model.pb -> ../../../frozen_model.pb
|-- in.lammps
|-- POSCAR -> ../relaxation/relax_task/CONTCAR
|-- task.000000
| | -- conf.lmp
| | -- frozen_model.pb -> ../frozen_model.pb
| | -- in.lammps -> ../in.lammps
| | -- inter.json
| | -- POSCAR
| | -- strain.json
|-- task.json
|-- task.000001
| | -- conf.lmp
| | -- frozen_model.pb -> ../frozen_model.pb
| | -- in.lammps -> ../in.lammps
| | -- inter.json
| | -- POSCAR
| | -- strain.json
|-- task.json
...
|-- task.000023
| -- conf.lmp
| -- frozen_model.pb -> ../frozen_model.pb
| -- in.lammps -> ../in.lammps
| -- inter.json
| -- POSCAR
| -- strain.json
|-- task.json
```

equi.stress.json records the stress information of the equilibrium task and strain.json records the deformation information of the corresponding task.

#### Vacancy output:

```
confs/std-fcc/vacancy_00/
```

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```

|-- frozen_model.pb -> ../../../../frozen_model.pb
|-- in.lammps
|-- POSCAR -> ../relaxation/relax_task/CONTCAR
|-- task.000000
|   |-- conf.lmp
|   |-- frozen_model.pb -> ../frozen_model.pb
|   |-- in.lammps -> ../in.lammps
|   |-- inter.json
|   |-- POSCAR
|   |-- supercell.json
|-- task.json

```

supercell.json records the supercell size information of the corresponding task.

#### Interstitial output:

```

confs/std-fcc/interstitial_00/
|-- element.out
|-- frozen_model.pb -> ../../../../frozen_model.pb
|-- in.lammps
|-- POSCAR -> ../relaxation/relax_task/CONTCAR
|-- task.000000
|   |-- conf.lmp
|   |-- frozen_model.pb -> ../frozen_model.pb
|   |-- in.lammps -> ../in.lammps
|   |-- inter.json
|   |-- POSCAR
|   |-- supercell.json
|-- task.json
|-- task.000001
|   |-- conf.lmp
|   |-- frozen_model.pb -> ../frozen_model.pb
|   |-- in.lammps -> ../in.lammps
|   |-- inter.json
|   |-- POSCAR
|   |-- supercell.json
|-- task.json

```

element.out records the inserted element type of each task and supercell.json records the supercell size information of the corresponding task.

#### Surface output:

```

confs/std-fcc/surface_00/
|-- frozen_model.pb -> ../../../../frozen_model.pb
|-- in.lammps
|-- POSCAR -> ../relaxation/relax_task/CONTCAR
|-- task.000000
|   |-- conf.lmp
|   |-- frozen_model.pb -> ../frozen_model.pb
|   |-- in.lammps -> ../in.lammps
|   |-- inter.json
|   |-- miller.json

```

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```

| | -- POSCAR
| | -- POSCAR.tmp
| | -- task.json
|-- task.000001
| | -- conf.lmp
| | -- frozen_model.pb -> ../frozen_model.pb
| | -- in.lammps -> ../in.lammps
| | -- inter.json
| | -- miller.json
| | -- POSCAR
| | -- POSCAR.tmp
| | -- task.json
...
|-- task.000008
| | -- conf.lmp
| | -- frozen_model.pb -> ../frozen_model.pb
| | -- in.lammps -> ../in.lammps
| | -- inter.json
| | -- miller.json
| | -- POSCAR
| | -- POSCAR.tmp
| | -- task.json

```

miller.json records the miller index of the corresponding task.

### 7.4.3 Property run

```
nohup dpngen autotest run property.json machine-ali.json > run.result 2>&1 &
```

the result file log.lammps, dump.relax, and outlog would be sent back.

### 7.4.4 Property-post

Use command

```
dpngen autotest post property.json
```

to post results as result.json and result.out in each property's path.

### 7.4.5 Properties

#### EOS get started and input examples

Equation of State (EOS) here calculates the energies of the most stable structures as a function of volume. Users can refer to Figure 4 of the [dpngen CPC paper](#) for more information of EOS.

### An example of the input file for EOS by VASP:

```
{
  "structures":      ["confs/mp-*", "confs/std-*", "confs/test-*"],
  "interaction": {
    "type":          "vasp",
    "incar":         "vasp_input/INCAR",
    "potcar_prefix": "vasp_input",
    "potcars":       {"Al": "POTCAR.al", "Mg": "POTCAR.mg"}
  },
  "properties": [
    {
      "type":         "eos",
      "vol_start":    0.9,
      "vol_end":      1.1,
      "vol_step":     0.01
    }
  ]
}
```

vol\_start is the starting volume relative to the equilibrium structure, vol\_step is the volume increment step relative to the equilibrium structure, and the biggest relative volume is smaller than vol\_end.

### EOS make

**Step 1.** Before make in EOS, the equilibrium configuration CONTCAR must be present in confs/mp-\*/relaxation.

**Step 2.** For the input example in the previous section, when we do make, 40 tasks would be generated as confs/mp-\*/eos\_00/task.000000, confs/mp-\*/eos\_00/task.000001, ... , confs/mp-\*/eos\_00/task.000039. The suffix 00 is used for possible refine later.

**Step 3.** If the task directory, for example confs/mp-\*/eos\_00/task.000000 is not empty, the old input files in it including INCAR, POSCAR, POTCAR, conf.lmp, in.lammps would be deleted.

**Step 4.** In each task directory, POSCAR.orig would link to confs/mp-\*/relaxation/CONTCAR. Then the scale parameter can be calculated as:

```
scale = (vol_current / vol_equi) ** (1. / 3.)
```

vol\_current is the corresponding volume per atom of the current task and vol\_equi is the volume per atom of the equilibrium configuration. Then the poscar\_scale function in dpgen.auto\_test.lib.vasp module would help to generate POSCAR file with vol\_current in confs/mp-\*/eos\_00/task.[0-9]\*[0-9].

**Step 5.** According to the task type, the input file including INCAR, POTCAR or conf.lmp, in.lammps would be written in every confs/mp-\*/eos\_00/task.[0-9]\*[0-9].

## EOS run

The work path of each task should be in the form like `confs/mp-*/eos_00` and all task is in the form like `confs/mp-*/eos_00/task.[0-9]*[0-9]`.

When we dispatch tasks, we would go through every individual work path in the list `confs/mp-*/eos_00`, and then submit `task.[0-9]*[0-9]` in each work path.

## EOS post

The post processing of EOS would go to every directory in `confs/mp-*/eos_00` and do the post processing. Let's suppose we are now in `confs/mp-100/eos_00` and there are `task.000000`, `task.000001`, ..., `task.000039` in this directory. By reading `inter.json` file in every task directory, the task type can be determined and the energy and force information of every task can further be obtained. By appending the dict of energy and force into a list, an example of the list with 1 atom is given as:

```
[
  {"energy": E1, "force": [fx1, fy1, fz1]},
  {"energy": E2, "force": [fx2, fy2, fz2]},
  ...
  {"energy": E40, "force": [fx40, fy40, fz40]}
]
```

Then the volume can be calculated from the task id and the corresponding energy can be obtained from the list above. Finally, there would be `result.json` in json format and `result.out` in txt format in `confs/mp-100/eos_00` containing the EOS results.

An example of `result.json` is give as:

```
{
  "14.808453313267595": -3.7194474,
  "14.972991683415014": -3.7242038,
  ...
  "17.934682346068534": -3.7087655
}
```

An example of `result.out` is given below:

```
onf_dir: /root/auto_test_example/deepmd/confs/std-fcc/eos_00
VpA(A^3)  EpA(eV)
14.808    -3.7194
14.973    -3.7242
...
17.935    -3.7088
```



## Elastic get started and input examples

Here we calculate the mechanical properties which include elastic constants (C11 to C66), bulk modulus  $B_v$ , shear modulus  $G_v$ , Youngs modulus  $E_v$ , and Poission ratio  $U_v$  of a certain crystal structure.

### An example of the input file for Elastic by deepmd:

```
{
  "structures": ["confs/mp-*", "confs/std-*", "confs/test-*"],
  "interaction": {
    "type": "deepmd",
    "model": "frozen_model.pb",
    "type_map": {"Al": 0, "Mg": 1}
  },
  "properties": [
    {
      "type": "elastic",
      "norm_deform": 1e-2,
      "shear_deform": 1e-2
    }
  ]
}
```

Here the default values of `norm_deform` and `shear_deform` are **1e-2** and **1e-2**, respectively. A list of `norm_strains` and `shear_strains` would be generated as below:

```
[-norm_def, -0.5 * norm_def, 0.5 * norm_def, norm_def]
[-shear_def, -0.5 * shear_def, 0.5 * shear_def, shear_def]
```

## Elastic make

**Step 1.** The `DeformedStructureSet` module in `pymatgen.analysis.elasticity.strain` is used to generate a set of independently deformed structures. `equi.stress.out` file is written to record the equilibrium stress in the Elastic directory. For the example in the previous section, `equi.stress.out` should be in `confs/mp-*/elastic_00`.

**Step 2.** If there are `init_from_suffix` and `output_suffix` parameter in the `properties` part, the *refine process* follows. Else, the deformed structure (POSCAR) and strain information (`strain.out`) are written in the task directory, for example, in `confs/mp-*/elastic_00/task.0000000`.

**Step 3.** When doing elastic by VASP, `ISIF=2`. When doing by LAMMPS, the following `in.lammps` would be written.

```
units          metal
dimension      3
boundary       p      p      p
atom_style     atomic
box            tilt large
read_data      conf.lmp
mass           1 1
mass           2 1
neigh_modify   every 1 delay 0 check no
pair_style     deepmd frozen_model.pb
```

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```

pair_coeff
compute      mype all pe
thermo        100
thermo_style  custom step pe pxx pyy pzz pxy pxz pyz lx ly lz vol c_mype
dump          1 all custom 100 dump.relax id type xs ys zs fx fy fz
min_style     cg
minimize      0 1.0000000e-10 5000 500000
variable      N equal count(all)
variable      V equal vol
variable      E equal "c_mype"
variable      Pxx equal pxx
variable      Pyy equal pyy
variable      Pzz equal pzz
variable      Pxy equal pxy
variable      Pxz equal pxz
variable      Pyz equal pyz
variable      Epa equal ${E}/${N}
variable      Vpa equal ${V}/${N}
print "All done"
print "Total number of atoms = ${N}"
print "Final energy per atoms = ${Epa}"
print "Final volume per atoms = ${Vpa}"
print "Final Stress (xx yy zz xy xz yz) = ${Pxx} ${Pyy} ${Pzz} ${Pxy} ${Pxz} ${Pyz}"

```

## Elastic run

Very similar to the run operation of EOS except for in different directories. Now the work path of each task should be in the form like `confs/mp-*/elastic_00` and all task is in the form like `confs/mp-*/elastic_00/task.[0-9]*[0-9]`.

## Elastic post

The ElasticTensor module in `pymatgen.analysis.elasticity.elastic` is used to get the elastic tensor, Bv, and Gv. The mechanical properties of a crystal structure would be written in `result.json` in json format and `result.out` in txt format. The example of the output file is give below.

## result.json

```

{
  "elastic_tensor": [
    134.90955999999997,
    54.329958699999985,
    51.802386099999985,
    3.5745279599999993,
    -1.38863259999999648e-05,
    -1.96382339999999486e-05,
    54.558402999999999,
    134.59654699999996,
    51.7972336,

```

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```

-3.53972684,
1.8395687999999963e-05,
8.7567993999999951e-05,
51.91324859999999,
51.913292199999994,
137.01763799999998,
-5.0903393999999969e-05,
6.992516299999996e-05,
3.7364786999999946e-05,
3.8780564440000007,
-3.770445632,
-1.2766205999999956,
35.41343199999999,
2.2479590800000023e-05,
1.3837692000000017e-06,
-4.959999999495933e-06,
2.5800000003918792e-06,
1.4800000030874965e-06,
2.9000000008417968e-06,
35.375960199999994,
3.8608356,
0.0,
0.0,
0.0,
0.0,
4.02554856,
38.375018399999995
],
"BV": 80.315363022222,
"GV": 38.40582656,
"EV": 99.37716395728943,
"uV": 0.2937771799031088
}

```

The order of `elastic_tensor` is C11, C12, ..., C16, C21, C22, ..., C26, ..., C66 and the unit of Bv, Gv, Ev, and uv is GPa.

## result.out

```

/root/auto_test_example/deepmd/confs/std-fcc/elastic_00
134.91  54.33  51.80  3.57  -0.00  -0.00
54.56  134.60  51.80  -3.54  0.00  0.00
51.91  51.91  137.02  -0.00  0.00  0.00
3.88   -3.77  -1.28  35.41  0.00  0.00
-0.00  0.00  0.00  0.00  35.38  3.86
0.00   0.00  0.00  0.00  4.03  38.38
# Bulk   Modulus BV = 80.32 GPa
# Shear  Modulus GV = 38.41 GPa
# Youngs Modulus EV = 99.38 GPa
# Poission Ratio uV = 0.29

```

## Vacancy get started and input examples

Vacancy calculates the energy difference when removing an atom from the crystal structure. We only need to give the information of `supercell` to help calculate the vacancy energy and the default value of `supercell` is `[1, 1, 1]`.

### An example of the input file for Vacancy by deepmd:

```
{
  "structures":      "confs/mp-*",
  "interaction": {
    "type":          "deepmd",
    "model":          "frozen_model.pb",
    "type_map":       {"Al": 0, "Mg": 1}
  },
  "properties": [
    {
      "type":         "vacancy",
      "supercell":     [1, 1, 1]
    }
  ]
}
```

## Vacancy make

**Step 1.** The `VacancyGenerator` module in `pymatgen.analysis.defects.generators` is used to generate a set of structures with vacancy.

**Step 2.** If there are `init_from_suffix` and `output_suffix` parameter in the `properties` part, the *refine process* follows. If `reproduce` is evoked, the *reproduce process* follows. Otherwise, the vacancy structure (POSCAR) and supercell information (`supercell.out`) are written in the task directory, for example, in `confs/mp-*/vacancy_00/task.000000` with the check and possible removing of the old input files like before.

**Step 3.** When doing vacancy by VASP, `ISIF = 3`. When doing vacancy by LAMMPS, the same `in.lammps` as that in *EOS* (*change\_box is True*) would be generated with `scale` set to one.

## Vacancy run

Very similar to the run operation of EOS except for in different directories. Now the work path of each task should be in the form like `confs/mp-*/vacancy_00` and all task is in the form like `confs/mp-*/vacancy_00/task.[0-9]*[0-9]`.

## Vacancy post

For Vacancy, we need to calculate the energy difference between a crystal structure with and without a vacancy. The examples of the output files `result.json` in json format and `result.out` in txt format are given below.

**result.json**

```
{
  "[3, 3, 3]-task.000000": [
    0.7352769999999964,
    -96.644642,
    -97.379919
  ]
}
```

**result.out**

```
/root/auto_test_example/deepmd/confs/std-fcc/vacancy_00
Structure:      Vac_E(eV)  E(eV)  equi_E(eV)
[3, 3, 3]-task.000000:  0.735  -96.645  -97.380
```

**Interstitial get started and input examples**

Interstitial calculates the energy difference when adding an atom into the crystal structure. We need to give the information of supercell (default value is [1, 1, 1]) and insert\_ele list for the element types of the atoms added in.

**An example of the input file for Interstitial by deepmd:**

```
{
  "structures":      "confs/mp-*",
  "interaction": {
    "type":          "deepmd",
    "model":          "frozen_model.pb",
    "type_map":       {"Al": 0, "Mg": 1}
  },
  "properties": [
    {
      "type":         "interstitial",
      "supercell":    [3, 3, 3],
      "insert_ele":   ["Al"],
      "conf_filters": {"min_dist": 1.5},
      "cal_setting":  {"input_prop": "lammmps_input/lammmps_high"}
    }
  ]
}
```

We add a `conf_filters` parameter in `properties` part and this parameter can help to eliminate undesirable structure which can render rather difficult convergence in calculations. In the example above, “**min\_dist**”: 1.5 means if the smallest atomic distance in the structure is less than 1.5 angstrom, the configuration would be eliminated and not used in calculations.

## Interstitial make

**Step 1.** For each element in `insert_ele` list, `InterstitialGenerator` module in `pymatgen.analysis.defects.generators` would help to generate interstitial structure. The structure would be appended into a list if it can meet the requirements in `conf_filters`.

**Step 2.** If `refine` is True, we do *refine process*. If `reprod-opt` is True (the default is **False**), we do *reproduce process*. Else, the vacancy structure (POSCAR) and supercell information (`supercell.out`) are written in the task directory, for example, in `confs/mp-*/interstitial_00/task.0000000` with the check and possible removing of the old input files like before.

**Step 3.** In interstitial by VASP, `ISIF = 3`. In interstitial by LAMMPS, the same `in.lammps` as that in *EOS (change\_box is True)* would be generated with `scale` set to one.

## Interstitial run

Very similar to the run operation of EOS except for in different directories. Now the work path of each task should be in the form like `confs/mp-*/interstitial_00` and all task is in the form like `confs/mp-*/interstitial_00/task.[0-9]*[0-9]`.

## Interstitial post

For Interstitial, we need to calculate the energy difference between a crystal structure with and without atom added in. The examples of the output files `result.json` in json format and `result.out` in txt format are given below.

### result.json

```
{
  "Al-[3, 3, 3]-task.0000000": [
    4.0229520000000004,
    -100.84773,
    -104.870682
  ],
  "Al-[3, 3, 3]-task.0000001": [
    2.78295200000000088,
    -102.08773,
    -104.870682
  ]
}
```

### result.out

```
/root/auto_test_example/deepmd/confs/std-fcc/interstitial_00
Insert_ele-Struct: Inter_E(eV)  E(eV)  equi_E(eV)
Al-[3, 3, 3]-task.0000000:   4.023  -100.848 -104.871
Al-[3, 3, 3]-task.0000001:   2.783  -102.088 -104.871
```

## Surface get started and input examples

Surface calculates the surface energy. We need to give the information of `min_slab_size`, `min_vacuum_size`, `max_miller` (default value is 2), and `pert_xz` which means perturbations in xz and will help work around vasp bug.

### An example of the input file for Surface by deepmd:

```
{
  "structures":      "confs/mp-*",
  "interaction": {
    "type":          "deepmd",
    "model":          "frozen_model.pb",
    "type_map":      {"Al": 0, "Mg": 1}
  },
  "properties": [
    {
      "type":         "surface",
      "min_slab_size": 10,
      "min_vacuum_size": 11,
      "max_miller":    2,
      "cal_type":      "static"
    }
  ]
}
```

## Surface make

**Step 1.** Based on the equilibrium configuration, `generate_all_slabs` module in `pymatgen.core.surface` would help to generate surface structure list with using `max_miller`, `min_slab_size`, and `min_vacuum_size` parameters.

**Step 2.** If `refine` is True, we do *refine process*. If `reprod-opt` is True (the default is False), we do *reproduce process*. Otherwise, the surface structure (POSCAR) with perturbations in xz and miller index information (`miller.out`) are written in the task directory, for example, in `confs/mp-*/interstitial_00/task.000000` with the check and possible removing of the old input files like before.

## Surface run

Very similar to the run operation of EOS except for in different directories. Now the work path of each task should be in the form like `confs/mp-*/surface_00` and all task is in the form like `confs/mp-*/surface_00/task.[0-9]*[0-9]`.

## Surface post

For Surface, we need to calculate the energy difference between a crystal structure with and without a surface with a certain miller index divided by the surface area.

The examples of the output files `result.json` in json format and `result.out` in txt format are given below.

## result.json

```
{
  "[1, 1, 1]-task.000000": [
    0.8051037974207992,
    -3.6035018,
    -3.7453815
  ],
  "[2, 2, 1]-task.000001": [
    0.9913881928811771,
    -3.5781115999999997,
    -3.7453815
  ],
  "[1, 1, 0]-task.000002": [
    0.9457333586026173,
    -3.5529366000000002,
    -3.7453815
  ],
  "[2, 2, -1]-task.000003": [
    0.9868013100872397,
    -3.5590607142857142,
    -3.7453815
  ],
  "[2, 1, 1]-task.000004": [
    1.0138239046484236,
    -3.563035875,
    -3.7453815
  ],
  "[2, 1, -1]-task.000005": [
    1.0661817319108005,
    -3.5432459166666668,
    -3.7453815
  ],
  "[2, 1, -2]-task.000006": [
    1.034003253044026,
    -3.550884125,
    -3.7453815
  ],
  "[2, 0, -1]-task.000007": [
```

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```

    0.9569958287615818,
    -3.5685403333333334,
    -3.7453815
  ],
  "[2, -1, -1]-task.000008": [
    0.9432935501134583,
    -3.5774615714285716,
    -3.7453815
  ]
}

```

## result.out

```

/root/auto_test_example/deepmd/confs/std-fcc/surface_00
Miller_Indices:      Surf_E(J/m^2)  EpA(eV)  equi_EpA(eV)
[1, 1, 1]-task.000000:      0.805      -3.604   -3.745
[2, 2, 1]-task.000001:      0.991      -3.578   -3.745
[1, 1, 0]-task.000002:      0.946      -3.553   -3.745
[2, 2, -1]-task.000003:      0.987      -3.559   -3.745
[2, 1, 1]-task.000004:      1.014      -3.563   -3.745
[2, 1, -1]-task.000005:      1.066      -3.543   -3.745
[2, 1, -2]-task.000006:      1.034      -3.551   -3.745
[2, 0, -1]-task.000007:      0.957      -3.569   -3.745
[2, -1, -1]-task.000008:      0.943      -3.577   -3.745

```

## 7.5 Refine

### 7.5.1 Refine get started and input examples

Sometimes we want to refine the calculation of a property from previous results. For example, when higher convergence criteria EDIFF and EDIFFG are necessary in VASP, the new VASP calculation is desired to start from the previous output configuration, rather than starting from scratch.

An example of the input file `refine.json` is given below:

```

{
  "structures":      ["confs/std-*"],
  "interaction": {
    "type":          "deepmd",
    "model":          "frozen_model.pb",
    "type_map":      {"Al": 0}
  },
  "properties": [
    {
      "type":          "vacancy",
      "init_from_suffix": "00",
      "output_suffix":  "01",
      "cal_setting":    {"input_prop": "lammps_input/lammps_high"}
    }
  ]
}

```

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```

    }
  ]
}

```

In this example, `refine` would output the results to `vacancy_01` based on the previous results in `vacancy_00` by using a different input commands file for `lammps`.

### 7.5.2 Refine make

```

dpgen autotest make refine.json
tree confs/std-fcc/vacancy_01/

```

the output will be:

```

confs/std-fcc/vacancy_01/
|-- frozen_model.pb -> ../../../../frozen_model.pb
|-- in.lammps
|-- task.000000
|   |-- conf.lmp
|   |-- frozen_model.pb -> ../frozen_model.pb
|   |-- in.lammps -> ../in.lammps
|   |-- inter.json
|   |-- POSCAR -> ../../vacancy_00/task.000000/CONTCAR
|   |-- supercell.json -> ../../vacancy_00/task.000000/supercell.json
|-- task.json

```

an new directory `vacancy_01` would be established and the starting configuration links to previous results.

### 7.5.3 Refine run

```

nohup dpgen autotest run refine.json machine-ali.json > run.result 2>&1 &

```

the run process of `refine` is similar to before.

### 7.5.4 Refine post

```

dpgen autotest post refine.json

```

the post process of `refine` is similar to the corresponding property.

## 7.6 Reproduce

### 7.6.1 Reproduce get started and input examples

Sometimes we want to reproduce the initial results with the same configurations for cross validation. This version of autotest package can accomplish this successfully in all property types except for Elastic. An input example for using deepmd to reproduce the VASP Interstitial results is given below:

```
{
  "structures":      ["confs/std-*"],
  "interaction": {
    "type":          "deepmd",
    "model":          "frozen_model.pb",
    "type_map":       {"Al": 0}
  },
  "properties": [
    {
      "type":          "interstitial",
      "reproduce":       true,
      "init_from_suffix": "00",
      "init_data_path": "../vasp/confs",
      "reprod_last_frame": false
    }
  ]
}
```

reproduce denotes whether to do reproduce or not and the default value is False.

init\_data\_path is the path of VASP or LAMMPS initial data to be reproduced. init\_from\_suffix is the suffix of the initial data and the default value is "00". In this case, the VASP Interstitial results are stored in ../vasp/confs/std-\*/interstitial\_00 and the reproduced Interstitial results would be in deepmd/confs/std-\*/interstitial\_reprod.

reprod\_last\_frame denotes if only the last frame is used in reproduce. The default value is True for eos and surface, but is False for vacancy and interstitial.

### 7.6.2 Reproduce make

```
dpngen autotest make reproduce.json
tree confs/std-fcc/interstitial_reprod/
```

the output will be:

```
confs/std-fcc/interstitial_reprod/
|-- frozen_model.pb -> ../../../../frozen_model.pb
|-- in.lammps
|-- task.0000000
|  |-- conf.lmp
|  |-- frozen_model.pb -> ../frozen_model.pb
|  |-- in.lammps -> ../in.lammps
|  |-- inter.json
|  |-- POSCAR
```

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```

|  -- task.json
|-- task.000001
|  -- conf.lmp
|  -- frozen_model.pb -> ../frozen_model.pb
|  -- in.lammps -> ../in.lammps
|  -- inter.json
|  -- POSCAR
|  -- task.json
...
|-- task.000038
|  -- conf.lmp
|  -- frozen_model.pb -> ../frozen_model.pb
|  -- in.lammps -> ../in.lammps
|  -- inter.json
|  -- POSCAR
|-- task.json

```

every single frame in the initial data is split into each task and the following `in.lammps` would help to do the static calculation:

```

clear
units          metal
dimension      3
boundary        p p p
atom_style      atomic
box            tilt large
read_data       conf.lmp
mass           1 26.982
neigh_modify    every 1 delay 0 check no
pair_style      deepmd frozen_model.pb
pair_coeff
compute         mype all pe
thermo          100
thermo_style    custom step pe pxx pyy pzz pxy pxz pyz lx ly lz vol c_mype
dump            1 all custom 100 dump.relax id type xs ys zs fx fy fz
run             0
variable        N equal count(all)
variable        V equal vol
variable        E equal "c_mype"
variable        tmp1x equal lx
variable        tmp1y equal ly
variable        Pxx equal pxx
variable        Pyy equal pyy
variable        Pzz equal pzz
variable        Pxy equal pxy
variable        Pxz equal pxz
variable        Pyz equal pyz
variable        Epa equal ${E}/${N}
variable        Vpa equal ${V}/${N}
variable        AA equal (${tmp1x}*${tmp1y})
print "All done"
print "Total number of atoms = ${N}"

```

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```

print "Final energy per atoms = ${Epa}"
print "Final volume per atoms = ${Vpa}"
print "Final Base area = ${AA}"
print "Final Stress (xx yy zz xy xz yz) = ${Pxx} ${Pyy} ${Pzz} ${Pxy} ${Pxz} ${Pyz}"

```

### 7.6.3 Reproduce run

```
nohup dpngen autotest run reproduce.json machine-ali.json > run.result 2>&1 &
```

the run process of reproduce is similar to before.

### 7.6.4 Reproduce post

```
dpngen autotest post reproduce.json
```

the output will be:

result.out:

```

/root/auto_test_example/deepmd/confs/std-fcc/interstitial_reprod
Reproduce: Initial_path Init_E(eV/atom)  Reprod_E(eV/atom)  Difference(eV/atom)
.../vasp/confs/std-fcc/interstitial_00/task.000000 -3.020 -3.240 -0.220
.../vasp/confs/std-fcc/interstitial_00/task.000000 -3.539 -3.541 -0.002
.../vasp/confs/std-fcc/interstitial_00/task.000000 -3.582 -3.582 -0.001
.../vasp/confs/std-fcc/interstitial_00/task.000000 -3.582 -3.581 0.001
.../vasp/confs/std-fcc/interstitial_00/task.000000 -3.594 -3.593 0.001
.../vasp/confs/std-fcc/interstitial_00/task.000000 -3.594 -3.594 0.001
.../vasp/confs/std-fcc/interstitial_00/task.000000 -3.598 -3.597 0.001
.../vasp/confs/std-fcc/interstitial_00/task.000000 -3.600 -3.600 0.001
.../vasp/confs/std-fcc/interstitial_00/task.000000 -3.600 -3.600 0.001
.../vasp/confs/std-fcc/interstitial_00/task.000000 -3.601 -3.600 0.001
.../vasp/confs/std-fcc/interstitial_00/task.000000 -3.602 -3.601 0.001
.../vasp/confs/std-fcc/interstitial_00/task.000000 -3.603 -3.602 0.001
.../vasp/confs/std-fcc/interstitial_00/task.000000 -3.603 -3.602 0.001
.../vasp/confs/std-fcc/interstitial_00/task.000000 -3.603 -3.602 0.001
.../vasp/confs/std-fcc/interstitial_00/task.000000 -3.603 -3.602 0.001
.../vasp/confs/std-fcc/interstitial_00/task.000000 -3.603 -3.602 0.001
.../vasp/confs/std-fcc/interstitial_00/task.000000 -3.603 -3.602 0.001
.../vasp/confs/std-fcc/interstitial_00/task.000001 -3.345 -3.372 -0.027
.../vasp/confs/std-fcc/interstitial_00/task.000001 -3.546 -3.556 -0.009
.../vasp/confs/std-fcc/interstitial_00/task.000001 -3.587 -3.593 -0.007
.../vasp/confs/std-fcc/interstitial_00/task.000001 -3.593 -3.599 -0.006
.../vasp/confs/std-fcc/interstitial_00/task.000001 -3.600 -3.606 -0.006
.../vasp/confs/std-fcc/interstitial_00/task.000001 -3.600 -3.606 -0.006
.../vasp/confs/std-fcc/interstitial_00/task.000001 -3.624 -3.631 -0.006
.../vasp/confs/std-fcc/interstitial_00/task.000001 -3.634 -3.640 -0.007
.../vasp/confs/std-fcc/interstitial_00/task.000001 -3.637 -3.644 -0.007
.../vasp/confs/std-fcc/interstitial_00/task.000001 -3.637 -3.644 -0.007
.../vasp/confs/std-fcc/interstitial_00/task.000001 -3.638 -3.645 -0.007

```

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```

.../vasp/confs/std-fcc/interstitial_00/task.000001 -3.638 -3.645 -0.007
.../vasp/confs/std-fcc/interstitial_00/task.000001 -3.639 -3.646 -0.007
.../vasp/confs/std-fcc/interstitial_00/task.000001 -3.639 -3.646 -0.007
.../vasp/confs/std-fcc/interstitial_00/task.000001 -3.639 -3.646 -0.007
.../vasp/confs/std-fcc/interstitial_00/task.000001 -3.639 -3.646 -0.007
.../vasp/confs/std-fcc/interstitial_00/task.000001 -3.639 -3.646 -0.007
.../vasp/confs/std-fcc/interstitial_00/task.000001 -3.639 -3.646 -0.007
.../vasp/confs/std-fcc/interstitial_00/task.000001 -3.639 -3.646 -0.007
.../vasp/confs/std-fcc/interstitial_00/task.000001 -3.639 -3.646 -0.007
.../vasp/confs/std-fcc/interstitial_00/task.000001 -3.639 -3.646 -0.007

```

the comparison of the initial and reproduced results as well as the absolute path of the initial data is recorded.

result.json:

```

{
  "/root/auto_test_example/vasp/confs/std-fcc/interstitial_00/task.000000": {
    "nframes": 18,
    "error": 0.0009738182472213228
  },
  "/root/auto_test_example/vasp/confs/std-fcc/interstitial_00/task.000001": {
    "nframes": 21,
    "error": 0.0006417039154057605
  }
}

```

the error analysis corresponding to the initial data is recorded and the error of the first frame is disregarded when all the frames are considered in reproduce.

## USER GUIDE

This part aims to show you how to get the community's help. Some frequently asked questions are listed in troubleshooting, and the explanation of errors that often occur is listed in common errors. If other unexpected problems occur, you're welcome to contact us for help.

### 8.1 Discussions

Welcome everyone to participate in the discussion about DP-GEN in the [discussion](#) module. You can ask for help, share an idea or anything to discuss here. Note: before you raise a question, please check TUTORIAL/FAQs and search history discussions to find solutions.

### 8.2 Issue

If you want to make a bug report or a request for new features, you can make an issue in the issue module.

Here are the types you can choose. A proper type can help developer figure out what you need. Also, you can assign yourself to solve the issue. Your contribution is welcome!

Note: before you raise a question, please check TUTORIAL/FAQs and search history issues to find solutions.

### 8.3 Tutorials

Tutorials can be found [here](#).

### 8.4 Example for parameters

If you have no idea how to prepare a PARAM for your task, you can find examples of PARAM for different tasks in [examples](#).

For example, if you want to set specific template for LAMMPS, you can find an example [here](#)

If you want to learn more about Machine parameters, please check [docs for dpdispatcher](#)

## 8.5 Pull requests - How to contribute

## 8.6 Troubleshooting

1. The most common problem is whether two settings correspond with each other, including:
  - The order of elements in `type_map` and `mass_map` and **`fp_pp_files`**.
  - Size of `init_data_sys` and `init_batch_size`.
  - Size of `sys_configs` and `sys_batch_size`.
  - Size of `sel_a` and actual types of atoms in your system.
  - Index of `sys_configs` and `sys_idx`.
2. Please verify the directories of `sys_configs`. If there isn't any POSCAR for `01.model_devi` in one iteration, it may happen that you write the false path of `sys_configs`. Note that `init_data_sys` is a list, while `sys_configs` should be a two-dimensional list. The first dimension corresponds to `sys_idx`, and the second level are some poscars under each group. Refer to the [sample file](#).
3. Correct format of JSON file.
4. The frames of one system should be larger than `batch_size` and `numb_test` in `default_training_param`. It happens that one iteration adds only a few structures and causes error in next iteration's training. In this condition, you may let `fp_task_min` be larger than `numb_test`.
5. If you found the dpgen with the same version on two machines behaves differently, you may have modified the code in one of them.

## 8.7 Common Errors

(Errors are sorted alphabetically)

### 8.7.1 Command not found: xxx.

There is no such software in the environment, or it is unavailable. It may be because 1. It is not installed; 2. The Conda environment is not activated; 3. You have chosen the wrong image in machine.json.

### 8.7.2 `dargs.dargs.ArgumentKeyError: [at location xxx] undefined key xxx is not allowed in strict mode.`

Strict format check has been applied since version 0.10.7. To avoid misleading users, some older-version keys that are already ignored or absorbed into default settings are not allowed to be present. And the expected structure of the dictionary in the param.json also differs from those before version 0.10.7. This error will occur when format check finds older-fashion keys in the json file. Please try deleting or annotating these keys, or correspondingly modulate the json file. Example files in the newest format could be found in [examples](#).



### 8.7.3 dargs.dargs.ArgumentTypeError: [at root location] key xxx gets wrong value type, requires but gets

Please check your parameters with [DPGEN's Document](#). Maybe you have superfluous parentheses in your parameter file.

### 8.7.4 FileNotFoundError: [Errno 2] No such file or directory: '.../01.model\_devi/graph.xxx.pb'

If you find this error occurs, please check your initial data. Your model will not be generated if the initial data is incorrect.

### 8.7.5 json.decoder.JSONDecodeError

Your .json file is incorrect. It may be a mistake in syntax or a missing comma.

### 8.7.6 OSError: [Error cannot find valid a data system] Please check your setting for data systems

Check if the path to the dataset in the parameter file is set correctly. Note that `init_data_sys` is a list, while `sys_configs` should be a two-dimensional list. The first dimension corresponds to `sys_idx`, and the second level are some poscars under each group. Refer to the [sample file](#).

### 8.7.7 RuntimeError: job:xxxxxxx failed 3 times

```
RuntimeError: job:xxxxxxx failed 3 times

.....

RuntimeError: Meet errors will handle unexpected submission state.
Debug information: remote_root==xxxxxxx
Debug information: submission_hash==xxxxxxx
Please check the dirs and scripts in remote_root. The job information mentioned above,
↪may help.
```

If a user finds an error like this, he or she is advised to check the files on the remote server. It shows that your job has failed 3 times, but has not shown the reason.

To find the reason, you can check the log on the remote root. For example, you can check `train.log`, which is generated by DeepMD-kit. It can tell you more details. If it doesn't help, you can manually run the `.sub` script, whose path is shown in Debug information: `remote_root==xxxxxxx`

Some common reasons are as follows:

1. Two or more jobs are submitted manually or automatically at the same time, and their hash value collide. This bug will be fixed in `dpdispatcher`.
2. You may have something wrong in your input files, which causes the process to fail.

### **8.7.8 RuntimeError: find too many unsuccessfully terminated jobs.**

The ratio of failed jobs is larger than `ratio_failure`. You can set a high value for `ratio_failure` or check if there is something wrong with your input files.

### **8.7.9 ValueError: Cannot load file containing picked data when `allow_picked=False`**

Please ensure that you write the correct path of the dataset with no excess files.

### **8.7.10 warnings.warn("Some Gromacs commands were NOT found; "**

You can ignore this warning if you don't need Gromacs. It just show that Gromacs is not installed in you environment.

## CONTRIBUTING GUIDE

### 9.1 Contributing Guide

The way to make contributions is through making pull requests(PR for short). After your PR is merged, the changes you make can be applied by other users.

Firstly, fork in DP-GEN repository. Then you can clone the repository, build a new branch, make changes and then make a pull request.

---

#### 9.1.1 How to contribute to DP-GEN

Welcome to the repository of [DP-GEN](#)

DP-GEN adopts the same convention as other software in DeepModeling Community.

You can first refer to DeepPMD-kit's [Contributing guide](#) and [Developer guide](#).

You can also read relative chapters on [Github Docs](#).

If you have no idea how to fix your problem or where to find the relative source code, please check [Code Structure](#) of the DP-GEN repository on this website.

#### Use command line

You can use git with the command line, or open the repository on Github Desktop. Here is a video as a demo of making changes to DP-GEN and publishing it with command line.

If you have never used Github before, remember to generate your ssh key and configure the public key in Github Settings. If you can't configure your username and password, please use token. The explanation from Github see [Github Blog: token authentication requirements for git operations](#). A discussion on [StackOverflow](#) can solve this problem.

## Use Github Desktop

Also, you can use Github Desktop to make PR. The following shows the steps to clone the repository and add your doc to tutorials. If it is your first time using Github, Open with Github Desktop is recommended. Github Desktop is a software, which can make your operations on branches visually.

After you clone it to your PC, you can open it with Github Desktop.

Firstly, create your new branch based on devel branch.

Secondly, add your doc to the certain directory in your local repository, and add its name into index.

Here is an [example](#). Remember to add the filename of your doc into index!

Thirdly, select the changes that you want to push, and commit to it. Press “Publish branch” to push your origin repository to the remote branch.

Finally, you can check it on github and make a pull request. Press “Compare & pull request” to make a PR.

(Note: please commit pr to the devel branch)

## 9.1.2 How to contribute to DP-GEN tutorials and documents

Welcome to [the documents of DP-GEN](#)

- If you want to add the documentation of a toy model, simply put your file in the directory doc/toymodels/ and push;
- If you want to add a new directory for a new category of instructions, make a new directory and add it in doc/index.rst.

Also welcome to [Tutorials repository](#) You can find the structure of tutorials and preparations before writing a document in [Writing Tips](#).

The latest page of DP-GEN Docs

### Examples of contributions

- [Example 1](#)
- [Example 2](#) (a simple one for beginner)

#### 1. Push your doc

#### 2. Add the directory in index.rst

#### 3. Build and check it

As mentioned in “How to build the website to check if the modification works”.

## 4. Make pull request to dpngen

### 9.1.3 Find how a parameter is used in the code

It is strongly recommended that you use the `find in files` function of Visual Studio software, Search function of Visual Studio Code, or similar functions of other software. Type in the name of the parameter you are looking for, and you will see where it is read in and used in the procedure. Of course, you can also search for the relevant code according to the above guide.

### 9.1.4 Want to modify a function?

If you have special requirements, you can make personalized modifications in the code corresponding to the function. If you think your modification can benefit the public, and it does not conflict with the current DP-GEN function; or if you fix a bug, please make a pull request to contribute the optimization to the DP-GEN repository.

### 9.1.5 DP-GEN dependencies

`dpdispatcher` and `dpdata` are dependencies of DP-GEN. `dpdispatcher` is related to task submission, monitoring and recovery, and `dpdata` is related to data processing. If you encounter an error and want to find the reason, please judge whether the problem comes from DP-GEN, `dpdispatcher` or `dpdata` according to the last line of `Traceback`.

### 9.1.6 About the update of the parameter file

You may have noticed that there are `arginfo.py` files in many folders. This is a file used to generate parameter documentation. If you add or modify a parameter in DP-GEN and intend to export it to the main repository, please sync your changes in `arginfo`.

### 9.1.7 Tips

1. Please try to submit a PR after finishing all the changes
2. Please briefly describe what you do with `git commit -m "<conclude-the-change-you-make>"`! “No description provided.” will make the maintainer feel confused.
3. It is not recommended to make changes directly in the `devel` branch. It is recommended to pull a branch from `devel`: `git checkout -b <new-branch-name>`
4. When switching branches, remember to check if you want to bring the changes to the next branch!
5. Please fix the errors reported by the unit test. You can firstly test on your local machine before pushing commits. Hint: The way to test the code is to go from the main directory to the tests directory, and use the command `python3 -m unittest`. You can watch the demo video for review. Sometimes you may fail unit tests due to your local circumstance. You can check whether the error reported is related to the part you modified to eliminate this problem. After submitting, as long as there is a green check mark after the PR title on the webpage, it means that the test has been passed.
6. Pay attention to whether there are comments under your PR. If there is a change request, you need to check and modify the code. If there are conflicts, you need to solve them manually.

---

After successfully making a PR, developers will check it and give comments. It will be merged after everything done. Then CONGRATULATIONS! You become a first-time contributor to DP-GEN!

How to get help from the community



## 10.1 dpgen package

`dpgen.info()`

### 10.1.1 Subpackages

`dpgen.auto_test` package

Subpackages

`dpgen.auto_test.lib` package

Submodules

`dpgen.auto_test.lib.abacus` module

`dpgen.auto_test.lib.abacus.check_finished(fname)`

`dpgen.auto_test.lib.abacus.check_stru_fixed(struf, fixed)`

`dpgen.auto_test.lib.abacus.final_stru(abacus_path)`

`dpgen.auto_test.lib.abacus.make_kspacing_kpt(struf, kspacing)`

`dpgen.auto_test.lib.abacus.modify_stru_path(strucf, tpath)`

`dpgen.auto_test.lib.abacus.poscar2stru(poscar, inter_param, stru='STRU')`

- `poscar`: POSCAR for input
- `inter_param`: dictionary of 'interaction' from `param.json`  
some key words for ABACUS are:
  - `atom_masses`: a dictionary of atoms' masses
  - `orb_files`: a dictionary of orbital files
  - `deepks_desc`: a string of deepks descriptor file
- `stru`: output filename, usually is 'STRU'.

`dpgen.auto_test.lib.abacus.stru2Structure(struf)`

```
dpngen.auto_test.lib.abacus.stru_fix_atom(struf, fix_atom=[True, True, True])
... ATOMIC_POSITIONS Cartesian #Cartesian(Unit is LATTICE_CONSTANT) Si #Name of element 0.0
#Magnetic for this element. 2 #Number of atoms 0.00 0.00 0.00 0 0 0 #x,y,z, move_x, move_y, move_z 0.25
0.25 0.25 0 0 0.

dpngen.auto_test.lib.abacus.stru_scale(stru_in, stru_out, scale)

dpngen.auto_test.lib.abacus.write_input(inputf, inputdict)

dpngen.auto_test.lib.abacus.write_kpt(kptf, kptlist)
```

### dpngen.auto\_test.lib.crys module

```
dpngen.auto_test.lib.crys.bcc(ele_name='ele', a=3.2144871302356037)

dpngen.auto_test.lib.crys.dhcp(ele_name='ele', a=2.863782463805517, c=9.353074360871936)

dpngen.auto_test.lib.crys.diamond(ele_name='ele', a=2.551340126037118)

dpngen.auto_test.lib.crys.fcc(ele_name='ele', a=4.05)

dpngen.auto_test.lib.crys.fcc1(ele_name='ele', a=4.05)

dpngen.auto_test.lib.crys.hcp(ele_name='ele', a=2.863782463805517, c=4.676537180435968)

dpngen.auto_test.lib.crys.sc(ele_name='ele', a=2.551340126037118)
```

### dpngen.auto\_test.lib.lammps module

```
dpngen.auto_test.lib.lammps.apply_type_map(conf_file, deepmd_type_map, ptypes)
    Apply type map. conf_file: conf file converted from POSCAR deepmd_type_map: deepmd atom type map
    ptypes: atom types defined in POSCAR.

dpngen.auto_test.lib.lammps.check_finished(fname)

dpngen.auto_test.lib.lammps.check_finished_new(fname, keyword)

dpngen.auto_test.lib.lammps.cvt_lammps_conf(fin, fout, type_map, ofmt='lammps/data')
    Format convert from fin to fout, specify the output format by ofmt Imcomplete situation.

dpngen.auto_test.lib.lammps.element_list(type_map)

dpngen.auto_test.lib.lammps.get_base_area(log)
    Get base area.

dpngen.auto_test.lib.lammps.get_nev(log)
    Get natoms, energy_per_atom and volume_per_atom from lammps log.

dpngen.auto_test.lib.lammps.get_stress(log)
    Get stress from lammps log.

dpngen.auto_test.lib.lammps.inter_deepmd(param)

dpngen.auto_test.lib.lammps.inter_eam_alloy(param)

dpngen.auto_test.lib.lammps.inter_eam_fs(param)
```



```
dpngen.auto_test.lib.lammps.inter_meam(param)
```

```
dpngen.auto_test.lib.lammps.make_lammps_elastic(conf, type_map, interaction, param, etol=0,
                                                ftol=1e-10, maxiter=5000, maxeval=500000)
```

```
dpngen.auto_test.lib.lammps.make_lammps_equi(conf, type_map, interaction, param, etol=0, ftol=1e-10,
                                              maxiter=5000, maxeval=500000, change_box=True)
```

```
dpngen.auto_test.lib.lammps.make_lammps_eval(conf, type_map, interaction, param)
```

```
dpngen.auto_test.lib.lammps.make_lammps_phonon(conf, masses, interaction, param, etol=0, ftol=1e-10,
                                              maxiter=5000, maxeval=500000)
```

Make lammps input for elastic calculation.

```
dpngen.auto_test.lib.lammps.make_lammps_press_relax(conf, type_map, scale2equi, interaction, param,
                                                    B0=70, bp=0, etol=0, ftol=1e-10, maxiter=5000,
                                                    maxeval=500000)
```

```
dpngen.auto_test.lib.lammps.poscar_from_last_dump(dump, poscar_out, deepmd_type_map)
```

Get poscar from the last frame of a lammps MD traj (dump format).

### dpngen.auto\_test.lib.lmp module

```
dpngen.auto_test.lib.lmp.box2lmpbox(orig, box)
```

```
dpngen.auto_test.lib.lmp.from_system_data(system)
```

```
dpngen.auto_test.lib.lmp.get_atoms(lines)
```

```
dpngen.auto_test.lib.lmp.get_atype(lines)
```

```
dpngen.auto_test.lib.lmp.get_lmpbox(lines)
```

```
dpngen.auto_test.lib.lmp.get_natoms(lines)
```

```
dpngen.auto_test.lib.lmp.get_natoms_vec(lines)
```

```
dpngen.auto_test.lib.lmp.get_natomtypes(lines)
```

```
dpngen.auto_test.lib.lmp.get_posi(lines)
```

```
dpngen.auto_test.lib.lmp.lmpbox2box(lohi, tilt)
```

```
dpngen.auto_test.lib.lmp.system_data(lines)
```

```
dpngen.auto_test.lib.lmp.to_system_data(lines)
```

### dpngen.auto\_test.lib.mfp\_eosfit module

```
dpngen.auto_test.lib.mfp_eosfit.BM4(vol, pars)
```

Birch-Murnaghan 4 pars equation from PRB 70, 224107, 3-order.

```
dpngen.auto_test.lib.mfp_eosfit.BM5(vol, pars)
```

Birch-Murnaghan 5 pars equation from PRB 70, 224107, 4-Order.

`dpngen.auto_test.lib.mfp_eosfit.LOG4(vol, pars)`

Natrual strain (Poirier-Tarantola)EOS with 4 paramters Seems only work in near-equilibrium range.

`dpngen.auto_test.lib.mfp_eosfit.LOG5(vol, parameters)`

Natrual strain (Poirier-Tarantola)EOS with 5 paramters.

`dpngen.auto_test.lib.mfp_eosfit.Li4p(V, parameters)`

Li JH, APL, 87, 194111 (2005).

`dpngen.auto_test.lib.mfp_eosfit.SJX_5p(vol, par)`

SJX\_5p's five parameters EOS, Physica B: Condens Mater, 2011, 406: 1276-1282.

`dpngen.auto_test.lib.mfp_eosfit.SJX_v2(vol, par)`

Sun Jiuxun, et al. J phys Chem Solids, 2005, 66: 773-782. They said it is satisfied for the limiting condition at high pressure.

`dpngen.auto_test.lib.mfp_eosfit.TEOS(v, par)`

Holland, et al, Journal of Metamorphic Geology, 2011, 29(3): 333-383 Modified Tait equation of Huang & Chow.

`dpngen.auto_test.lib.mfp_eosfit.birch(v, parameters)`

From Intermetallic compounds: Principles and Practice, Vol. I: Principles Chapter 9 pages 195-210 by M. Mehl. B. Klein, D. Papaconstantopoulos paper downloaded from Web.

case where n=0

`dpngen.auto_test.lib.mfp_eosfit.calc_props_BM4(pars)`

`dpngen.auto_test.lib.mfp_eosfit.calc_props_LOG4(pars)`

`dpngen.auto_test.lib.mfp_eosfit.calc_props_SJX_5p(par)`

`dpngen.auto_test.lib.mfp_eosfit.calc_props_mBM4(pars)`

`dpngen.auto_test.lib.mfp_eosfit.calc_props_mBM4poly(pars)`

`dpngen.auto_test.lib.mfp_eosfit.calc_props_mBM5poly(pars)`

`dpngen.auto_test.lib.mfp_eosfit.calc_props_morse(pars)`

`dpngen.auto_test.lib.mfp_eosfit.calc_props_morse_6p(par)`

`dpngen.auto_test.lib.mfp_eosfit.calc_props_vinet(pars)`

`dpngen.auto_test.lib.mfp_eosfit.calc_v0_mBM4poly(x, pars)`

`dpngen.auto_test.lib.mfp_eosfit.calc_v0_mBM5poly(x, pars)`

`dpngen.auto_test.lib.mfp_eosfit.ext_splint(xp, yp, order=3, method='unispl')`

`dpngen.auto_test.lib.mfp_eosfit.ext_vec(func, fin, p0, fs, fe, vols=None, vole=None, ndata=101, refit=0, show_fig=False)`

Extrapolate the data points for E-V based on the fitted parameters in small or very large volume range.

`dpngen.auto_test.lib.mfp_eosfit.ext_velp(fin, fstart, fend, vols, vole, ndata, order=3, method='unispl', fout='ext_velp.dat', show_fig=False)`

Extrapolate the lattice parameters based on input data.

`dpngen.auto_test.lib.mfp_eosfit.get_eos_list()`

---

```

dpgen.auto_test.lib.mfp_eosfit.get_eos_list_3p()
dpgen.auto_test.lib.mfp_eosfit.get_eos_list_4p()
dpgen.auto_test.lib.mfp_eosfit.get_eos_list_5p()
dpgen.auto_test.lib.mfp_eosfit.get_eos_list_6p()
dpgen.auto_test.lib.mfp_eosfit.init_guess(fin)
dpgen.auto_test.lib.mfp_eosfit.lsqfit_eos(func, fin, par, fstart, fend, show_fig=False, fout='EoSfit.out',
                                         refit=-1)

dpgen.auto_test.lib.mfp_eosfit.mBM4(vol, pars)
    Birch-Murnaghan 4 pars equation from PRB 70, 224107, 3-order BM.
dpgen.auto_test.lib.mfp_eosfit.mBM4poly(vol, parameters)
    Modified BM5 EOS, Shang SL comput mater sci, 2010: 1040-1048, original expressions.
dpgen.auto_test.lib.mfp_eosfit.mBM5(vol, pars)
    Modified BM5 EOS, Shang SL comput mater sci, 2010: 1040-1048.
dpgen.auto_test.lib.mfp_eosfit.mBM5poly(vol, pars)
    Modified BM5 EOS, Shang SL comput mater sci, 2010: 1040-1048, original expressions.
dpgen.auto_test.lib.mfp_eosfit.mie(v, p)
    Mie model for song's FVT.
dpgen.auto_test.lib.mfp_eosfit.mie_simple(v, p)
    Mie_simple model for song's FVT.
dpgen.auto_test.lib.mfp_eosfit.morse(v, pars)
    Reproduce from ShunliShang's matlab script.
dpgen.auto_test.lib.mfp_eosfit.morse_3p(volume, p)
    morse_AB EOS formula from Song's FVT souces A= 0.5*B.
dpgen.auto_test.lib.mfp_eosfit.morse_6p(vol, par)
    Generalized Morse EOS proposed by Qin, see: Qin et al. Phys Rev B, 2008, 78, 214108. Qin et al. Phys Rev B,
    2008, 77, 220103(R).
dpgen.auto_test.lib.mfp_eosfit.morse_AB(volume, p)
    morse_AB EOS formula from Song's FVT souces.
dpgen.auto_test.lib.mfp_eosfit.murnaghan(vol, pars)
    Four-parameters murnaghan EOS. From PRB 28,5480 (1983).
dpgen.auto_test.lib.mfp_eosfit.parse_argument()
dpgen.auto_test.lib.mfp_eosfit.rBM4(vol, pars)
    Implementions as Alberto Otero-de-la-Roza, i.e. rBM4 is used here Comput Physics Comm, 2011, 182: 1708-
    1720.
dpgen.auto_test.lib.mfp_eosfit.rBM4_pv(vol, pars)
    Implementions as Alberto Otero-de-la-Roza, i.e. rBM4 is used here Comput Physics Comm, 2011, 182: 1708-
    1720 Fit for V-P relations.

```

`dpgen.auto_test.lib.mfp_eosfit.rBM5(vol, pars)`

Implementions as Alberto Otero-de-la-Roza, i.e. rBM5 is used here Comput Physics Comm, 2011, 182: 1708-1720.

`dpgen.auto_test.lib.mfp_eosfit.rBM5_pv(vol, pars)`

Implementions as Alberto Otero-de-la-Roza, i.e. rBM5 is used here Comput Physics Comm, 2011, 182: 1708-1720 Fit for V-P relations.

`dpgen.auto_test.lib.mfp_eosfit.rPT4(vol, pars)`

Natrual strain EOS with 4 paramters Seems only work in near-equilibrium range. Implementions as Alberto Otero-de-la-Roza, i.e. rPT4 is used here Comput Physics Comm, 2011, 182: 1708-1720, in their article, labeled as PT3 (3-order), however, we mention it as rPT4 for 4-parameters EOS.

`dpgen.auto_test.lib.mfp_eosfit.rPT4_pv(vol, pars)`

Natrual strain (Poirier-Tarantola)EOS with 4 paramters Seems only work in near-equilibrium range. Implementions as Alberto Otero-de-la-Roza, i.e. rPT4 is used here Comput Physics Comm, 2011, 182: 1708-1720, in their article, labeled as PT3 (3-order), however, we mention it as rPT4 for 4-parameters EOS.

`dpgen.auto_test.lib.mfp_eosfit.rPT5(vol, pars)`

Natrual strain EOS with 4 paramters Seems only work in near-equilibrium range. Implementions as Alberto Otero-de-la-Roza, i.e. rPT5 is used here Comput Physics Comm, 2011, 182: 1708-1720, in their article, labeled as PT3 (3-order), however, we mention it as rPT5 for 4-parameters EOS.

`dpgen.auto_test.lib.mfp_eosfit.rPT5_pv(vol, pars)`

Natrual strain (Poirier-Tarantola)EOS with 5 paramters Implementions as Alberto Otero-de-la-Roza, i.e. rPT5 is used here Comput Physics Comm, 2011, 182: 1708-1720, in their article, labeled as PT3 (3-order), however, we mention it as rPT5 for 4-parameters EOS.

`dpgen.auto_test.lib.mfp_eosfit.read_ve(fin)`

`dpgen.auto_test.lib.mfp_eosfit.read_velp(fin, fstart, fend)`

`dpgen.auto_test.lib.mfp_eosfit.read_vlp(fin, fstart, fend)`

`dpgen.auto_test.lib.mfp_eosfit.repro_ve(func, vol_i, p)`

`dpgen.auto_test.lib.mfp_eosfit.repro_vp(func, vol_i, pars)`

`dpgen.auto_test.lib.mfp_eosfit.res_BM4(pars, y, x)`

`dpgen.auto_test.lib.mfp_eosfit.res_BM5(pars, y, x)`

`dpgen.auto_test.lib.mfp_eosfit.res_LOG4(pars, y, x)`

`dpgen.auto_test.lib.mfp_eosfit.res_LOG5(pars, y, x)`

`dpgen.auto_test.lib.mfp_eosfit.res_Li4p(p, y, x)`

`dpgen.auto_test.lib.mfp_eosfit.res_SJX_5p(p, e, v)`

`dpgen.auto_test.lib.mfp_eosfit.res_SJX_v2(p, e, v)`

`dpgen.auto_test.lib.mfp_eosfit.res_TEOS(p, e, v)`

`dpgen.auto_test.lib.mfp_eosfit.res_birch(pars, y, x)`

`dpgen.auto_test.lib.mfp_eosfit.res_mBM4(pars, y, x)`

`dpgen.auto_test.lib.mfp_eosfit.res_mBM4poly(pars, y, x)`

```

dpngen.auto_test.lib.mfp_eosfit.res_mBM5(pars, y, x)
dpngen.auto_test.lib.mfp_eosfit.res_mBM5poly(pars, y, x)
dpngen.auto_test.lib.mfp_eosfit.res_mie(p, e, v)
dpngen.auto_test.lib.mfp_eosfit.res_mie_simple(p, e, v)
dpngen.auto_test.lib.mfp_eosfit.res_morse(p, en, volume)
dpngen.auto_test.lib.mfp_eosfit.res_morse_3p(p, en, volume)
dpngen.auto_test.lib.mfp_eosfit.res_morse_6p(p, en, volume)
dpngen.auto_test.lib.mfp_eosfit.res_morse_AB(p, en, volume)
dpngen.auto_test.lib.mfp_eosfit.res_murnaghan(pars, y, x)
dpngen.auto_test.lib.mfp_eosfit.res_rBM4(pars, y, x)
dpngen.auto_test.lib.mfp_eosfit.res_rBM4_pv(pars, y, x)
dpngen.auto_test.lib.mfp_eosfit.res_rBM5(pars, y, x)
dpngen.auto_test.lib.mfp_eosfit.res_rBM5_pv(pars, y, x)
dpngen.auto_test.lib.mfp_eosfit.res_rPT4(pars, y, x)
dpngen.auto_test.lib.mfp_eosfit.res_rPT4_pv(pars, y, x)
dpngen.auto_test.lib.mfp_eosfit.res_rPT5(pars, y, x)
dpngen.auto_test.lib.mfp_eosfit.res_rPT5_pv(pars, y, x)
dpngen.auto_test.lib.mfp_eosfit.res_universal(pars, y, x)
dpngen.auto_test.lib.mfp_eosfit.res_vinet(pars, y, x)
dpngen.auto_test.lib.mfp_eosfit.res_vinet_pv(pars, y, x)
dpngen.auto_test.lib.mfp_eosfit.universal(vol, parameters)
    Universal equation of state(Vinet P et al., J. Phys.: Condens. Matter 1, p1941 (1989)).
dpngen.auto_test.lib.mfp_eosfit.vinet(vol, pars)
    Vinet equation from PRB 70, 224107 Following, Shang Shunli et al., comput mater sci, 2010: 1040-1048,
    original expressions.
dpngen.auto_test.lib.mfp_eosfit.vinet_pv(vol, pars)

```

### **dpngen.auto\_test.lib.pwscf module**

```

dpngen.auto_test.lib.pwscf.make_pwscf_input(sys_data, fp_pp_files, fp_params)

```

**dpngen.auto\_test.lib.siesta module**

`dpngen.auto_test.lib.siesta.make_siesta_input(sys_data, fp_pp_files, fp_params)`

**dpngen.auto\_test.lib.util module**

`dpngen.auto_test.lib.util.collect_task(all_task, task_type)`

`dpngen.auto_test.lib.util.get_machine_info(mdata, task_type)`

`dpngen.auto_test.lib.util.insert_data(task, task_type, username, file_name)`

`dpngen.auto_test.lib.util.make_work_path(jdata, task, reprod_opt, static, user)`

`dpngen.auto_test.lib.util.voigt_to_stress(inpt)`

**dpngen.auto\_test.lib.utils module**

`dpngen.auto_test.lib.utils.cmd_append_log(cmd, log_file)`

`dpngen.auto_test.lib.utils.copy_file_list(file_list, from_path, to_path)`

`dpngen.auto_test.lib.utils.create_path(path)`

`dpngen.auto_test.lib.utils.log_iter(task, ii, jj)`

`dpngen.auto_test.lib.utils.log_task(message)`

`dpngen.auto_test.lib.utils.make_iter_name(iter_index)`

`dpngen.auto_test.lib.utils.record_iter(record, confs, ii, jj)`

`dpngen.auto_test.lib.utils.repeat_to_length(string_to_expand, length)`

`dpngen.auto_test.lib.utils.replace(file_name, pattern, subst)`

**dpngen.auto\_test.lib.vasp module**

**exception** `dpngen.auto_test.lib.vasp.OutcarItemError`

Bases: `Exception`

`dpngen.auto_test.lib.vasp.check_finished(fname)`

`dpngen.auto_test.lib.vasp.get_boxes(fname)`

`dpngen.auto_test.lib.vasp.get_energies(fname)`

`dpngen.auto_test.lib.vasp.get_nev(fname)`

`dpngen.auto_test.lib.vasp.get_poscar_natoms(fname)`

`dpngen.auto_test.lib.vasp.get_poscar_types(fname)`

`dpngen.auto_test.lib.vasp.get_stress(fname)`

```

dpgen.auto_test.lib.vasp.make_kspacing_kpoints(poscar, kspacing, kgamma)
dpgen.auto_test.lib.vasp.make_vasp_kpoints(kpoints, kgamma=False)
dpgen.auto_test.lib.vasp.make_vasp_kpoints_from_incar(work_dir, jdata)
dpgen.auto_test.lib.vasp.make_vasp_phonon_incar(ecut, ediff, npar, kpar, kspacing=0.5, kgamma=True,
                                                ismear=1, sigma=0.2)
dpgen.auto_test.lib.vasp.make_vasp_relax_incar(ecut, ediff, relax_ion, relax_shape, relax_volume,
                                                npar, kpar, kspacing=0.5, kgamma=True, ismear=1,
                                                sigma=0.22)
dpgen.auto_test.lib.vasp.make_vasp_static_incar(ecut, ediff, npar, kpar, kspacing=0.5, kgamma=True,
                                                ismear=1, sigma=0.2)
dpgen.auto_test.lib.vasp.perturb_xz(poscar_in, poscar_out, pert=0.01)
dpgen.auto_test.lib.vasp.poscar_natoms(poscar_in)
dpgen.auto_test.lib.vasp.poscar_scale(poscar_in, poscar_out, scale)
dpgen.auto_test.lib.vasp.poscar_vol(poscar_in)
dpgen.auto_test.lib.vasp.reciprocal_box(box)
dpgen.auto_test.lib.vasp.regulate_poscar(poscar_in, poscar_out)
dpgen.auto_test.lib.vasp.sort_poscar(poscar_in, poscar_out, new_names)

```

## Submodules

### dpgen.auto\_test.ABACUS module

```
class dpgen.auto_test.ABACUS.ABACUS(inter_parameter, path_to_poscar)
```

Bases: [Task](#)

## Methods

<a href="#"><i>backward_files</i></a> ([property_type])	Return backward files.
<a href="#"><i>compute</i></a> (output_dir)	Compute output of the task.
<a href="#"><i>forward_common_files</i></a> ([property_type])	Return forward common files.
<a href="#"><i>forward_files</i></a> ([property_type])	Return forward files.
<a href="#"><i>make_input_file</i></a> (output_dir, task_type, ...)	Prepare input files for a computational task For example, the VASP prepares INCAR.
<a href="#"><i>make_potential_files</i></a> (output_dir)	Prepare potential files for a computational task.

### modify\_input

```
backward_files(property_type='relaxation')
```

Return backward files.

**compute**(*output\_dir*)

Compute output of the task. IMPORTANT: The output configuration should be converted and stored in a CONTCAR file.

**Parameters****output\_dir**

[str] The directory storing the input and output files.

**Returns****result\_dict: dict**

A dict that storing the result. For example: { "energy": xxx, "force": [xxx] }

**Notes**

The following files are generated: CONTCAR: output file

The output configuration is converted to CONTCAR and stored in the *output\_dir*

**forward\_common\_files**(*property\_type='relaxation'*)

Return forward common files.

**forward\_files**(*property\_type='relaxation'*)

Return forward files.

**make\_input\_file**(*output\_dir, task\_type, task\_param*)

Prepare input files for a computational task For example, the VASP prepares INCAR. LAMMPS (including DeePMD, MEAM...) prepares in.lammps.

**Parameters****output\_dir**

[str] The directory storing the input files.

**task\_type**

[str] Can be - "relaxation": structure relaxation - "static": static computation calculates the energy, force... of a strcture

**task\_param**

[dict] The parameters of the task. For example the VASP interaction can be provided with { "ediff": 1e-6, "ediffg": 1e-5 }

**make\_potential\_files**(*output\_dir*)

Prepare potential files for a computational task. For example, the VASP prepares POTCAR. DeePMD prepares frozen model(s). IMPORTANT: Interaction should be stored in *output\_dir/inter.json*.

**Parameters****output\_dir**

[str] The directory storing the potential files.

**Notes**

The following files are generated:

**inter.json: output file**

The task information is stored in *output\_dir/inter.json*

**modify\_input**(*incard, x, y*)



## dpngen.auto\_test.EOS module

**class** dpngen.auto\_test.EOS.**EOS**(parameter, inter\_param=None)

Bases: *Property*

### Methods

<code>compute(output_file, print_file, path_to_work)</code>	Postprocess the finished tasks to compute the property.
<code>make_confs(path_to_work, path_to_equi[, refine])</code>	Make configurations needed to compute the property.
<code>post_process(task_list)</code>	post_process the KPOINTS file in elastic.
<code>task_param()</code>	Return the parameter of each computational task, for example, {'ediffg': 1e-4}.
<code>task_type()</code>	Return the type of each computational task, for example, 'relaxation', 'static'....

**make\_confs**(path\_to\_work, path\_to\_equi, refine=False)

Make configurations needed to compute the property. The tasks directory will be named as path\_to\_work/task.xxxxxx IMPORTANT: handel the case when the directory exists.

#### Parameters

##### **path\_to\_work**

[str] The path where the tasks for the property are located

##### **path\_to\_equi**

[str] -refine == False: The path to the directory that equilibrated the configuration.  
-refine == True: The path to the directory that has property confs.

##### **refine**

[str] To refine existing property confs or generate property confs from a equilibrated conf

#### Returns

##### **task\_list: list of str**

The list of task directories.

**post\_process**(task\_list)

post\_process the KPOINTS file in elastic.

**task\_param**()

Return the parameter of each computational task, for example, {'ediffg': 1e-4}.

**task\_type**()

Return the type of each computational task, for example, 'relaxation', 'static'....

## dpngen.auto\_test.Elastic module

**class** dpngen.auto\_test.Elastic.**Elastic**(parameter, inter\_param=None)

Bases: *Property*

## Methods

<code>compute(output_file, print_file, path_to_work)</code>	Postprocess the finished tasks to compute the property.
<code>make_confs(path_to_work, path_to_equi[, refine])</code>	Make configurations needed to compute the property.
<code>post_process(task_list)</code>	post_process the KPOINTS file in elastic.
<code>task_param()</code>	Return the parameter of each computational task, for example, {'ediffg': 1e-4}.
<code>task_type()</code>	Return the type of each computational task, for example, 'relaxation', 'static'....

**make\_confs**(*path\_to\_work*, *path\_to\_equi*, *refine=False*)

Make configurations needed to compute the property. The tasks directory will be named as *path\_to\_work/task.xxxxxx* IMPORTANT: handel the case when the directory exists.

### Parameters

#### **path\_to\_work**

[str] The path where the tasks for the property are located

#### **path\_to\_equi**

[str] -refine == False: The path to the directory that equilibrated the configuration.  
-refine == True: The path to the directory that has property confs.

#### **refine**

[str] To refine existing property confs or generate property confs from a equilibrated conf

### Returns

#### **task\_list: list of str**

The list of task directories.

**post\_process**(*task\_list*)

post\_process the KPOINTS file in elastic.

**task\_param**()

Return the parameter of each computational task, for example, {'ediffg': 1e-4}.

**task\_type**()

Return the type of each computational task, for example, 'relaxation', 'static'....

**dpgen.auto\_test.Gamma module**

**class** dpgen.auto\_test.Gamma.**Gamma**(*parameter*, *inter\_param=None*)

Bases: *Property*

Calculation of common gamma lines for bcc and fcc.

**Methods**

<code>compute(output_file, print_file, path_to_work)</code>	Postprocess the finished tasks to compute the property.
<code>make_confs(path_to_work, path_to_equi[, refine])</code>	Make configurations needed to compute the property.
<code>post_process(task_list)</code>	post_process the KPOINTS file in elastic.
<code>task_param()</code>	Return the parameter of each computational task, for example, {'ediffg': 1e-4}.
<code>task_type()</code>	Return the type of each computational task, for example, 'relaxation', 'static'....

<b>centralize_slab</b> <b>return_direction</b>
---

**static** `centralize_slab(slab)` → None

**make\_confs**(*path\_to\_work*, *path\_to\_equi*, *refine=False*)

Make configurations needed to compute the property. The tasks directory will be named as `path_to_work/task.xxxxxx` IMPORTANT: handel the case when the directory exists.

**Parameters**

**path\_to\_work**

[str] The path where the tasks for the property are located

**path\_to\_equi**

[str] -refine == False: The path to the directory that equilibrated the configuration.  
-refine == True: The path to the directory that has property confs.

**refine**

[str] To refine existing property confs or generate property confs from a equilibrated conf

**Returns**

**task\_list: list of str**

The list of task directories.

**post\_process**(*task\_list*)

post\_process the KPOINTS file in elastic.

**return\_direction**()

**task\_param**()

Return the parameter of each computational task, for example, {'ediffg': 1e-4}.

**task\_type**()

Return the type of each computational task, for example, 'relaxation', 'static'....

**dpngen.auto\_test.Interstitial module**

```
class dpngen.auto_test.Interstitial.Interstitial(parameter, inter_param=None)
```

Bases: *Property*

**Methods**

<code>compute(output_file, print_file, path_to_work)</code>	Postprocess the finished tasks to compute the property.
<code>make_confs(path_to_work, path_to_equi[, refine])</code>	Make configurations needed to compute the property.
<code>post_process(task_list)</code>	post_process the KPOINTS file in elastic.
<code>task_param()</code>	Return the parameter of each computational task, for example, {'ediffg': 1e-4}.
<code>task_type()</code>	Return the type of each computational task, for example, 'relaxation', 'static'....

```
make_confs(path_to_work, path_to_equi, refine=False)
```

Make configurations needed to compute the property. The tasks directory will be named as path\_to\_work/task.xxxxxx IMPORTANT: handel the case when the directory exists.

**Parameters****path\_to\_work**

[str] The path where the tasks for the property are located

**path\_to\_equi**

[str] -refine == False: The path to the directory that equilibrated the configuration.  
-refine == True: The path to the directory that has property confs.

**refine**

[str] To refine existing property confs or generate property confs from a equilibrated conf

**Returns****task\_list: list of str**

The list of task directories.

```
post_process(task_list)
```

post\_process the KPOINTS file in elastic.

```
task_param()
```

Return the parameter of each computational task, for example, {'ediffg': 1e-4}.

```
task_type()
```

Return the type of each computational task, for example, 'relaxation', 'static'....

**dpngen.auto\_test.Lammps module**

```
class dpngen.auto_test.Lammps.Lammps(inter_parameter, path_to_poscar)
```

Bases: *Task*

## Methods

<code>backward_files([property_type])</code>	Return backward files.
<code>compute(output_dir)</code>	Compute output of the task.
<code>forward_common_files([property_type])</code>	Return forward common files.
<code>forward_files([property_type])</code>	Return forward files.
<code>make_input_file(output_dir, task_type, ...)</code>	Prepare input files for a computational task For example, the VASP prepares INCAR.
<code>make_potential_files(output_dir)</code>	Prepare potential files for a computational task.

<b>set_inter_type_func</b> <b>set_model_param</b>
--

**backward\_files**(*property\_type*='relaxation')

Return backward files.

**compute**(*output\_dir*)

Compute output of the task. IMPORTANT: The output configuration should be converted and stored in a CONTCAR file.

**Parameters**

**output\_dir**

[str] The directory storing the input and output files.

**Returns**

**result\_dict: dict**

A dict that storing the result. For example: { "energy": xxx, "force": [xxx] }

## Notes

The following files are generated: CONTCAR: output file

The output configuration is converted to CONTCAR and stored in the *output\_dir*

**forward\_common\_files**(*property\_type*='relaxation')

Return forward common files.

**forward\_files**(*property\_type*='relaxation')

Return forward files.

**make\_input\_file**(*output\_dir, task\_type, task\_param*)

Prepare input files for a computational task For example, the VASP prepares INCAR. LAMMPS (including DeePMD, MEAM...) prepares in.lammps.

**Parameters**

**output\_dir**

[str] The directory storing the input files.

**task\_type**

[str] Can be - "relaxation": structure relaxation - "static": static computation calculates the energy, force... of a structure

**task\_param**

[dict] The parameters of the task. For example the VASP interaction can be provided with { "ediff": 1e-6, "ediffg": 1e-5 }

**make\_potential\_files**(*output\_dir*)

Prepare potential files for a computational task. For example, the VASP prepares POTCAR. DeePMD prepares frozen model(s). IMPORTANT: Interaction should be stored in *output\_dir/inter.json*.

**Parameters****output\_dir**

[str] The directory storing the potential files.

**Notes**

The following files are generated:

**inter.json: output file**

The task information is stored in *output\_dir/inter.json*

**set\_inter\_type\_func()****set\_model\_param()****dpgen.auto\_test.Property module****class dpgen.auto\_test.Property.Property**(*parameter*)

Bases: ABC

**Attributes****task\_param**

Return the parameter of each computational task, for example, {'ediffg': 1e-4}.

**task\_type**

Return the type of each computational task, for example, 'relaxation', 'static'....

**Methods**

<i>compute</i> ( <i>output_file</i> , <i>print_file</i> , <i>path_to_work</i> )	Postprocess the finished tasks to compute the property.
<i>make_confs</i> ( <i>path_to_work</i> , <i>path_to_equi</i> [, <i>refine</i> ])	Make configurations needed to compute the property.
<i>post_process</i> ( <i>task_list</i> )	post_process the KPOINTS file in elastic.

**compute**(*output\_file*, *print\_file*, *path\_to\_work*)

Postprocess the finished tasks to compute the property. Output the result to a json database.

**Parameters****output\_file:**

The file to output the property in json format

**print\_file:**

The file to output the property in txt format

**path\_to\_work:**

The working directory where the computational tasks locate.

**abstract make\_confs**(*path\_to\_work*, *path\_to\_equi*, *refine=False*)

Make configurations needed to compute the property. The tasks directory will be named as *path\_to\_work/task.xxxxxx* IMPORTANT: handel the case when the directory exists.

**Parameters****path\_to\_work**

[str] The path where the tasks for the property are located

**path\_to\_equi**

[str] -refine == False: The path to the directory that equilibrated the configuration.  
 -refine == True: The path to the directory that has property confs.

**refine**

[str] To refine existing property confs or generate property confs from a equilibrated conf

**Returns****task\_list: list of str**

The list of task directories.

**abstract post\_process(task\_list)**

post\_process the KPOINTS file in elastic.

**abstract property task\_param**

Return the parameter of each computational task, for example, {'ediffg': 1e-4}.

**abstract property task\_type**

Return the type of each computational task, for example, 'relaxation', 'static'....

**dpngen.auto\_test.Surface module**

**class** dpngen.auto\_test.Surface.**Surface**(parameter, inter\_param=None)

Bases: *Property*

**Methods**

<code>compute(output_file, print_file, path_to_work)</code>	Postprocess the finished tasks to compute the property.
<code>make_confs(path_to_work, path_to_equi[, refine])</code>	Make configurations needed to compute the property.
<code>post_process(task_list)</code>	post_process the KPOINTS file in elastic.
<code>task_param()</code>	Return the parameter of each computational task, for example, {'ediffg': 1e-4}.
<code>task_type()</code>	Return the type of each computational task, for example, 'relaxation', 'static'....

**make\_confs**(path\_to\_work, path\_to\_equi, refine=False)

Make configurations needed to compute the property. The tasks directory will be named as path\_to\_work/task.xxxxxx IMPORTANT: handel the case when the directory exists.

**Parameters****path\_to\_work**

[str] The path where the tasks for the property are located

**path\_to\_equi**

[str] -refine == False: The path to the directory that equilibrated the configuration.  
 -refine == True: The path to the directory that has property confs.

**refine**

[str] To refine existing property confs or generate property confs from a equilibrated conf

**Returns****task\_list: list of str**

The list of task directories.

**post\_process**(*task\_list*)

post\_process the KPOINTS file in elastic.

**task\_param**()

Return the parameter of each computational task, for example, {'ediffg': 1e-4}.

**task\_type**()

Return the type of each computational task, for example, 'relaxation', 'static'...

## dpgen.auto\_test.Task module

**class** dpgen.auto\_test.Task.**Task**(*inter\_parameter*, *path\_to\_poscar*)

Bases: ABC

### Attributes

***backward\_files***

Return backward files.

***forward\_common\_files***

Return forward common files.

***forward\_files***

Return forward files.

### Methods

<b><i>compute</i></b> ( <i>output_dir</i> )	Compute output of the task.
<b><i>make_input_file</i></b> ( <i>output_dir</i> , <i>task_type</i> , ...)	Prepare input files for a computational task For example, the VASP prepares INCAR.
<b><i>make_potential_files</i></b> ( <i>output_dir</i> )	Prepare potential files for a computational task.

**abstract property backward\_files**

Return backward files.

**abstract compute**(*output\_dir*)

Compute output of the task. IMPORTANT: The output configuration should be converted and stored in a CONTCAR file.

### Parameters

**output\_dir**

[str] The directory storing the input and output files.

### Returns

**result\_dict: dict**

A dict that storing the result. For example: { "energy": xxx, "force": [xxx] }



## Notes

The following files are generated: CONTCAR: output file

The output configuration is converted to CONTCAR and stored in the *output\_dir*

### **abstract property forward\_common\_files**

Return forward common files.

### **abstract property forward\_files**

Return forward files.

### **abstract make\_input\_file**(*output\_dir*, *task\_type*, *task\_param*)

Prepare input files for a computational task For example, the VASP prepares INCAR. LAMMPS (including DeePMD, MEAM...) prepares in.lammps.

#### **Parameters**

##### **output\_dir**

[str] The directory storing the input files.

##### **task\_type**

[str] Can be - “relaxation”: structure relaxation - “static”: static computation calculates the energy, force... of a strcture

##### **task\_param**

[dict] The parameters of the task. For example the VASP interaction can be provided with { “ediff”: 1e-6, “ediffg”: 1e-5 }

### **abstract make\_potential\_files**(*output\_dir*)

Prepare potential files for a computational task. For example, the VASP prepares POTCAR. DeePMD prepares frozen model(s). IMPORTANT: Interaction should be stored in *output\_dir/inter.json*.

#### **Parameters**

##### **output\_dir**

[str] The directory storing the potential files.

## Notes

The following files are generated:

### **inter.json: output file**

The task information is stored in *output\_dir/inter.json*

## dpngen.auto\_test.VASP module

**class** dpngen.auto\_test.VASP.VASP(*inter\_parameter*, *path\_to\_poscar*)

Bases: *Task*

## Methods

<i>backward_files</i> ([ <i>property_type</i> ])	Return backward files.
<i>compute</i> ( <i>output_dir</i> )	Compute output of the task.
<i>forward_common_files</i> ([ <i>property_type</i> ])	Return forward common files.
<i>forward_files</i> ([ <i>property_type</i> ])	Return forward files.
<i>make_input_file</i> ( <i>output_dir</i> , <i>task_type</i> , ...)	Prepare input files for a computational task For example, the VASP prepares INCAR.
<i>make_potential_files</i> ( <i>output_dir</i> )	Prepare potential files for a computational task.

**backward\_files**(*property\_type='relaxation'*)

Return backward files.

**compute**(*output\_dir*)

Compute output of the task. IMPORTANT: The output configuration should be converted and stored in a CONTCAR file.

**Parameters**

**output\_dir**

[str] The directory storing the input and output files.

**Returns**

**result\_dict: dict**

A dict that storing the result. For example: { "energy": xxx, "force": [xxx] }

**Notes**

The following files are generated: CONTCAR: output file

The output configuration is converted to CONTCAR and stored in the *output\_dir*

**forward\_common\_files**(*property\_type='relaxation'*)

Return forward common files.

**forward\_files**(*property\_type='relaxation'*)

Return forward files.

**make\_input\_file**(*output\_dir, task\_type, task\_param*)

Prepare input files for a computational task For example, the VASP prepares INCAR. LAMMPS (including DeePMD, MEAM...) prepares in.lammps.

**Parameters**

**output\_dir**

[str] The directory storing the input files.

**task\_type**

[str] Can be - "relaxation": structure relaxation - "static": static computation calculates the energy, force... of a strcture

**task\_param**

[dict] The parameters of the task. For example the VASP interaction can be provided with { "ediff": 1e-6, "ediffg": 1e-5 }

**make\_potential\_files**(*output\_dir*)

Prepare potential files for a computational task. For example, the VASP prepares POTCAR. DeePMD prepares frozen model(s). IMPORTANT: Interaction should be stored in *output\_dir/inter.json*.

**Parameters**

**output\_dir**

[str] The directory storing the potential files.

**Notes**

The following files are generated:

**inter.json: output file**

The task information is stored in *output\_dir/inter.json*

## dpngen.auto\_test.Vacancy module

**class** dpngen.auto\_test.Vacancy.**Vacancy**(parameter, inter\_param=None)

Bases: *Property*

### Methods

<code>compute(output_file, print_file, path_to_work)</code>	Postprocess the finished tasks to compute the property.
<code>make_confs(path_to_work, path_to_equi[, refine])</code>	Make configurations needed to compute the property.
<code>post_process(task_list)</code>	post_process the KPOINTS file in elastic.
<code>task_param()</code>	Return the parameter of each computational task, for example, {'ediffg': 1e-4}.
<code>task_type()</code>	Return the type of each computational task, for example, 'relaxation', 'static'....

**make\_confs**(path\_to\_work, path\_to\_equi, refine=False)

Make configurations needed to compute the property. The tasks directory will be named as path\_to\_work/task.xxxxxx IMPORTANT: handel the case when the directory exists.

#### Parameters

##### **path\_to\_work**

[str] The path where the tasks for the property are located

##### **path\_to\_equi**

[str] -refine == False: The path to the directory that equilibrated the configuration.  
-refine == True: The path to the directory that has property confs.

##### **refine**

[str] To refine existing property confs or generate property confs from a equilibrated conf

#### Returns

##### **task\_list: list of str**

The list of task directories.

**post\_process**(task\_list)

post\_process the KPOINTS file in elastic.

**task\_param**()

Return the parameter of each computational task, for example, {'ediffg': 1e-4}.

**task\_type**()

Return the type of each computational task, for example, 'relaxation', 'static'....

## dpngen.auto\_test.calculator module

dpngen.auto\_test.calculator.**make\_calculator**(inter\_parameter, path\_to\_poscar)

Make an instance of Task.

### dpgen.auto\_test.common\_equi module

`dpgen.auto_test.common_equi.make_equi`(*confs*, *inter\_param*, *relax\_param*)

`dpgen.auto_test.common_equi.post_equi`(*confs*, *inter\_param*)

`dpgen.auto_test.common_equi.run_equi`(*confs*, *inter\_param*, *mdata*)

### dpgen.auto\_test.common\_prop module

`dpgen.auto_test.common_prop.make_property`(*confs*, *inter\_param*, *property\_list*)

`dpgen.auto_test.common_prop.make_property_instance`(*parameters*, *inter\_param*)

Make an instance of Property.

`dpgen.auto_test.common_prop.post_property`(*confs*, *inter\_param*, *property\_list*)

`dpgen.auto_test.common_prop.run_property`(*confs*, *inter\_param*, *property\_list*, *mdata*)

`dpgen.auto_test.common_prop.worker`(*work\_path*, *all\_task*, *forward\_common\_files*, *forward\_files*,  
*backward\_files*, *mdata*, *inter\_type*)

### dpgen.auto\_test.gen\_confs module

`dpgen.auto_test.gen_confs.gen_alloy`(*eles*, *key*)

`dpgen.auto_test.gen_confs.gen_ele_std`(*ele\_name*, *ctype*)

`dpgen.auto_test.gen_confs.gen_element`(*ele\_name*, *key*)

`dpgen.auto_test.gen_confs.gen_element_std`(*ele\_name*)

`dpgen.auto_test.gen_confs.make_path_mp`(*ii*)

`dpgen.auto_test.gen_confs.test_fit`(*struct*, *data*)

### dpgen.auto\_test.mpdb module

`dpgen.auto_test.mpdb.check_apikey`()

`dpgen.auto_test.mpdb.get_structure`(*mp\_id*)

### dpgen.auto\_test.refine module

`dpgen.auto_test.refine.make_refine`(*init\_from\_suffix*, *output\_suffix*, *path\_to\_work*)

### **dpngen.auto\_test.reproduce module**

`dpngen.auto_test.reproduce.make_repro`(*inter\_param*, *init\_data\_path*, *init\_from\_suffix*, *path\_to\_work*,  
*reprod\_last\_frame=True*)

`dpngen.auto_test.reproduce.post_repro`(*init\_data\_path*, *init\_from\_suffix*, *all\_tasks*, *ptr\_data*,  
*reprod\_last\_frame=True*)

### **dpngen.auto\_test.run module**

`dpngen.auto_test.run.gen_test`(*args*)

`dpngen.auto_test.run.run_task`(*step*, *param\_file*, *machine\_file=None*)

### **dpngen.collect package**

#### **Submodules**

#### **dpngen.collect.collect module**

`dpngen.collect.collect.collect_data`(*target\_folder*, *param\_file*, *output*, *verbose=True*, *shuffle=True*,  
*merge=True*)

`dpngen.collect.collect.gen_collect`(*args*)

### **dpngen.data package**

#### **Subpackages**

#### **dpngen.data.tools package**

#### **Submodules**

#### **dpngen.data.tools.bcc module**

`dpngen.data.tools.bcc.gen_box`()

`dpngen.data.tools.bcc.numb_atoms`()

`dpngen.data.tools.bcc.poscar_unit`(*latt*)

**dpngen.data.tools.cessp2force\_lin module**

```
dpngen.data.tools.cessp2force_lin.Parser()
dpngen.data.tools.cessp2force_lin.get_outcar_files(directory, recursive)
dpngen.data.tools.cessp2force_lin.process_outcar_file_v5_dev(outcars, data, numbers, types,
                                                             max_types, elements=None,
                                                             windex=None, fout='potfit.configs')
dpngen.data.tools.cessp2force_lin.scan_outcar_file(file_handle)
dpngen.data.tools.cessp2force_lin.uniq(seq)
```

**dpngen.data.tools.create\_random\_disturb module**

```
dpngen.data.tools.create_random_disturb.RandomDisturbParser()
dpngen.data.tools.create_random_disturb.create_disturbs_abacus_dev(fin, nfile, dmax=1.0,
                                                                    etmax=0.1, ofmt='abacus',
                                                                    dstyle='uniform',
                                                                    write_d=False, diag=0)
dpngen.data.tools.create_random_disturb.create_disturbs_ase(fin, nfile, dmax=1.0, ofmt='lmp',
                                                             dstyle='uniform', write_d=False)
dpngen.data.tools.create_random_disturb.create_disturbs_ase_dev(fin, nfile, dmax=1.0, etmax=0.1,
                                                                ofmt='lmp', dstyle='uniform',
                                                                write_d=False, diag=0)
dpngen.data.tools.create_random_disturb.create_disturbs_atomsk(fin, nfile, dmax=1.0, ofmt='lmp')
dpngen.data.tools.create_random_disturb.create_random_alloys(fin, alloy_dist, ifmt='vasp',
                                                             ofmt='vasp')
```

In fact, atomsk also gives us the convinient tool to do this.

```
dpngen.data.tools.create_random_disturb.gen_random_disturb(dmax, a, b, dstyle='uniform')
dpngen.data.tools.create_random_disturb.gen_random_emat(etmax, diag=0)
dpngen.data.tools.create_random_disturb.random_range(a, b, ndata=1)
```

**dpngen.data.tools.diamond module**

```
dpngen.data.tools.diamond.gen_box()
dpngen.data.tools.diamond.numb_atoms()
dpngen.data.tools.diamond.poscar_unit(latt)
```

**dpgen.data.tools.fcc module**

```
dpgen.data.tools.fcc.gen_box()
dpgen.data.tools.fcc.numb_atoms()
dpgen.data.tools.fcc.poscar_unit(latt)
```

**dpgen.data.tools.hcp module**

```
dpgen.data.tools.hcp.gen_box()
dpgen.data.tools.hcp.numb_atoms()
dpgen.data.tools.hcp.poscar_unit(latt)
```

**dpgen.data.tools.io\_lammps module**

ASE Atoms convert to LAMMPS configuration Some functions are adapted from ASE lammpsrun.py.

```
dpgen.data.tools.io_lammps.ase2lammpsdata(atoms, typeids=None, fout='out.lmp')
dpgen.data.tools.io_lammps.car2dir(v, Ainv)
    Cartesian to direct coordinates.
dpgen.data.tools.io_lammps.convert_cell(ase_cell)
    Convert a parallel piped (forming right hand basis) to lower triangular matrix LAMMPS can accept. This function
    transposes cell matrix so the bases are column vectors.
dpgen.data.tools.io_lammps.convert_forces(forces0, cell0, cell_new)
dpgen.data.tools.io_lammps.convert_positions(pos0, cell0, cell_new, direct=False)
dpgen.data.tools.io_lammps.convert_stress(s6_0, cell0, cell_new)
dpgen.data.tools.io_lammps.dir2car(v, A)
    Direct to cartesian coordinates.
dpgen.data.tools.io_lammps.get_atoms_ntypes(atoms)
dpgen.data.tools.io_lammps.get_typeid(typeids, csymbol)
dpgen.data.tools.io_lammps.is_upper_triangular(mat)
    Test if 3x3 matrix is upper triangular LAMMPS has a rule for cell matrix definition.
dpgen.data.tools.io_lammps.set_atoms_typeids(atoms)
dpgen.data.tools.io_lammps.set_atoms_typeids_with_atomic_numbers(atoms)
dpgen.data.tools.io_lammps.stress6_to_stress9(s6)
dpgen.data.tools.io_lammps.stress9_to_stress6(s9)
```

**dpgen.data.tools.ovito\_file\_convert module**

**dpgen.data.tools.poscar\_copy module**

**dpgen.data.tools.sc module**

`dpgen.data.tools.sc.gen_box()`

`dpgen.data.tools.sc.numb_atoms()`

`dpgen.data.tools.sc.poscar_unit(latt)`

## Submodules

**dpgen.data.arginfo module**

`dpgen.data.arginfo.init_bulk_abacus_args()` → list[Argument]

`dpgen.data.arginfo.init_bulk_jdata_arginfo()` → Argument

Generate arginfo for dpgen init\_bulk jdata.

### Returns

**Argument**

dpgen init\_bulk jdata arginfo

`dpgen.data.arginfo.init_bulk_mdata_arginfo()` → Argument

Generate arginfo for dpgen init\_bulk mdata.

### Returns

**Argument**

arginfo

`dpgen.data.arginfo.init_bulk_variant_type_args()` → list[Variant]

`dpgen.data.arginfo.init_bulk_vasp_args()` → list[Argument]

`dpgen.data.arginfo.init_reaction_jdata_arginfo()` → Argument

Generate arginfo for dpgen init\_reaction jdata.

### Returns

**Argument**

dpgen init\_reaction jdata arginfo

`dpgen.data.arginfo.init_reaction_mdata_arginfo()` → Argument

Generate arginfo for dpgen init\_reaction mdata.

### Returns

**Argument**

arginfo

`dpgen.data.arginfo.init_surf_jdata_arginfo()` → Argument

Generate arginfo for dpgen init\_surf jdata.

### Returns

**Argument**

dpgen init\_surf jdata arginfo



`dpgen.data.arginfo.init_surf_mdata_arginfo()` → Argument

Generate arginfo for dpgen init\_surf mdata.

**Returns**

**Argument**

arginfo

## dpgen.data.gen module

`dpgen.data.gen.class_cell_type(jdata)`

`dpgen.data.gen.coll_abacus_md(jdata)`

`dpgen.data.gen.coll_vasp_md(jdata)`

`dpgen.data.gen.create_path(path, back=False)`

`dpgen.data.gen.gen_init_bulk(args)`

`dpgen.data.gen.make_abacus_md(jdata, mdata)`

`dpgen.data.gen.make_abacus_relax(jdata, mdata)`

`dpgen.data.gen.make_combines(dim, natoms)`

`dpgen.data.gen.make_scale(jdata)`

`dpgen.data.gen.make_scale_ABACUS(jdata)`

`dpgen.data.gen.make_super_cell(jdata)`

`dpgen.data.gen.make_super_cell_ABACUS(jdata, stru_data)`

`dpgen.data.gen.make_super_cell_STRU(jdata)`

`dpgen.data.gen.make_super_cell_poscar(jdata)`

`dpgen.data.gen.make_unit_cell(jdata)`

`dpgen.data.gen.make_unit_cell_ABACUS(jdata)`

`dpgen.data.gen.make_vasp_md(jdata, mdata)`

`dpgen.data.gen.make_vasp_relax(jdata, mdata)`

`dpgen.data.gen.out_dir_name(jdata)`

`dpgen.data.gen.pert_scaled(jdata)`

`dpgen.data.gen.place_element(jdata)`

`dpgen.data.gen.place_element_ABACUS(jdata, supercell_stru)`

`dpgen.data.gen.poscar_ele(poscar_in, poscar_out, eles, natoms)`

`dpgen.data.gen.poscar_natoms(lines)`

`dpgen.data.gen.poscar_scale(poscar_in, poscar_out, scale)`

```
dpngen.data.gen.poscar_scale_abacus(poscar_in, poscar_out, scale, jdata)
dpngen.data.gen.poscar_scale_cartesian(str_in, scale)
dpngen.data.gen.poscar_scale_direct(str_in, scale)
dpngen.data.gen.poscar_shuffle(poscar_in, poscar_out)
dpngen.data.gen.replace(file_name, pattern, subst)
dpngen.data.gen.run_abacus_md(jdata, mdata)
dpngen.data.gen.run_abacus_relax(jdata, mdata)
dpngen.data.gen.run_vasp_md(jdata, mdata)
dpngen.data.gen.run_vasp_relax(jdata, mdata)
dpngen.data.gen.shuffle_stru_data(supercell_stru)
dpngen.data.gen.stru_ele(supercell_stru, stru_out, eles, natoms, jdata, path_work)
```

### dpngen.data.reaction module

input: trajectory 00: ReaxFF MD (lammmps) 01: build dataset (mddatasetbuilder) 02: fp (gaussian) 03: convert to deepmd data output: data.

```
dpngen.data.reaction.convert_data(jdata)
dpngen.data.reaction.gen_init_reaction(args)
dpngen.data.reaction.link_fp_input()
dpngen.data.reaction.link_reaxff(jdata)
dpngen.data.reaction.link_trj(jdata)
    Link lammppstrj.
dpngen.data.reaction.make_lmp(jdata)
dpngen.data.reaction.run_build_dataset(jdata, mdata, log_file='build_log')
dpngen.data.reaction.run_fp(jdata, mdata, log_file='output', forward_common_files=[])
dpngen.data.reaction.run_reaxff(jdata, mdata, log_file='reaxff_log')
```

### dpngen.data.surf module

```
dpngen.data.surf.class_cell_type(jdata)
dpngen.data.surf.create_path(path)
dpngen.data.surf.gen_init_surf(args)
dpngen.data.surf.make_combines(dim, natoms)
dpngen.data.surf.make_scale(jdata)
```

```

dpgen.data.surf.make_super_cell_pymatgen(jdata)
dpgen.data.surf.make_unit_cell(jdata)
dpgen.data.surf.make_vasp_relax(jdata)
dpgen.data.surf.out_dir_name(jdata)
dpgen.data.surf.pert_scaled(jdata)
dpgen.data.surf.place_element(jdata)
dpgen.data.surf.poscar_ele(poscar_in, poscar_out, eles, natoms)
dpgen.data.surf.poscar_elong(poscar_in, poscar_out, elong, shift_center=True)
dpgen.data.surf.poscar_natoms(poscar_in)
dpgen.data.surf.poscar_scale(poscar_in, poscar_out, scale)
dpgen.data.surf.poscar_scale_cartesian(str_in, scale)
dpgen.data.surf.poscar_scale_direct(str_in, scale)
dpgen.data.surf.poscar_shuffle(poscar_in, poscar_out)
dpgen.data.surf.replace(file_name, pattern, subst)
dpgen.data.surf.run_vasp_relax(jdata, mdata)

```

### dpgen.database package

```

class dpgen.database.DPPotcar(symbols=None, functional='PBE', pp_file=None, pp_lists=None)
    Bases: MSONable

```

#### Methods

<a href="#"><code>as_dict()</code></a>	A JSON serializable dict representation of an object.
<a href="#"><code>from_dict(d)</code></a>	
<b>param d</b>	
Dict representation.	
<code>to_json()</code>	Returns a json string representation of the MSONable object.
<code>unsafe_hash()</code>	Returns an hash of the current object.
<code>validate_monty_v1(_MSONable__input_value)</code>	Pydantic validator with correct signature for pydantic v1.x
<code>validate_monty_v2(_MSONable__input_value, _)</code>	Pydantic validator with correct signature for pydantic v2.x

```

from_file
write_file

```

**as\_dict()**

A JSON serializable dict representation of an object.

**classmethod from\_dict(*d*)**

**Parameters**

**d** – Dict representation.

**Returns**

MSONable class.

**classmethod from\_file(*filename*)**

**write\_file(*filename*)**

**class** dpgen.database.**Entry**(*composition, calculator, inputs, data, entry\_id=None, attribute=None, tag=None*)

Bases: MSONable

An lightweight Entry object containing key computed data for storing purpose.

**Parameters**

**composition**

[Composition] Composition of the entry. For flexibility, this can take the form of all the typical input taken by a Composition, including a {symbol: amt} dict, a string formula, and others.

**inputs**

[dict] An dict of parameters associated with the entry. Defaults to None.

**data**

[dict] An dict of any additional data associated with the entry. Defaults to None.

**entry\_id**

[obj] An optional id to uniquely identify the entry.

**attribute**

Optional attribute of the entry. This can be used to specify that the entry is a newly found compound, or to specify a particular label for the entry, or else ... Used for further analysis and plotting purposes. An attribute can be anything but must be MSONable.

**Attributes**

**number\_element**

## Methods

<b>as_dict()</b>	A JSON serializable dict representation of an object.
<b>from_dict(<i>d</i>)</b>	
<b>param d</b>	
Dict representation.	
<b>to_json()</b>	Returns a json string representation of the MSONable object.
<b>unsafe_hash()</b>	Returns an hash of the current object.
<b>validate_monty_v1(_MSONable__input_value)</b>	Pydantic validator with correct signature for pydantic v1.x
<b>validate_monty_v2(_MSONable__input_value, _)</b>	Pydantic validator with correct signature for pydantic v2.x

**as\_dict()**

A JSON serializable dict representation of an object.

**classmethod** `from_dict(d)`

**Parameters**

**d** – Dict representation.

**Returns**

MSONable class.

**property** `number_element`

**class** `dpngen.database.VaspInput`(*incar, poscar, potcar, kpoints=None, optional\_files=None, \*\*kwargs*)

Bases: dict, MSONable

Class to contain a set of vasp input objects corresponding to a run.

**Args:**

`incar`: Incar object. `kpoints`: Kpoints object. `poscar`: Poscar object. `potcar`: Potcar object. `optional_files`: Other input files supplied as a dict of {  
    `filename`: object}. The object should follow standard pymatgen conventions in implementing a `as_dict()` and `from_dict` method.

## Methods

<code>as_dict()</code>	A JSON serializable dict representation of an object.
<code>clear()</code>	
<code>copy()</code>	
<code>from_dict(d)</code>	<b>param d</b> Dict representation.
<code>from_directory(input_dir[, optional_files])</code>	Read in a set of VASP input from a directory.
<code>fromkeys(iterable[, value])</code>	Create a new dictionary with keys from iterable and values set to value.
<code>get(key[, default])</code>	Return the value for key if key is in the dictionary, else default.
<code>items()</code>	
<code>keys()</code>	
<code>pop(key[, default])</code>	If the key is not found, return the default if given; otherwise, raise a KeyError.
<code>popitem(/)</code>	Remove and return a (key, value) pair as a 2-tuple.
<code>setdefault(key[, default])</code>	Insert key with a value of default if key is not in the dictionary.
<code>to_json()</code>	Returns a json string representation of the MSONable object.
<code>unsafe_hash()</code>	Returns an hash of the current object.
<code>update([E, ]**F)</code>	If E is present and has a .keys() method, then does: for k in E: D[k] = E[k] If E is present and lacks a .keys() method, then does: for k, v in E: D[k] = v In either case, this is followed by: for k in F: D[k] = F[k]
<code>validate_monty_v1(_MSONable__input_value)</code>	Pydantic validator with correct signature for pydantic v1.x
<code>validate_monty_v2(_MSONable__input_value, _)</code>	Pydantic validator with correct signature for pydantic v2.x
<code>values()</code>	
<code>write_input([output_dir, ...])</code>	Write VASP input to a directory.

**as\_dict()**

A JSON serializable dict representation of an object.

**classmethod from\_dict(d)****Parameters**

**d** – Dict representation.

**Returns**

MSONable class.

**static from\_directory(input\_dir, optional\_files=None)**

Read in a set of VASP input from a directory. Note that only the standard INCAR, POSCAR, POTCAR and KPOINTS files are read unless optional\_filenames is specified.

**Parameters**

**input\_dir**

[str] Directory to read VASP input from.

**optional\_files**

[dict] Optional files to read in as well as a dict of {filename: Object type}. Object type must have a static method from\_file.

**write\_input**(*output\_dir='.', make\_dir\_if\_not\_present=True*)

Write VASP input to a directory.

**Parameters****output\_dir**

[str] Directory to write to. Defaults to current directory (“.”).

**make\_dir\_if\_not\_present**

[bool] Create the directory if not present. Defaults to True.

**Submodules****dpngen.database.entry module****class** dpngen.database.entry.**Entry**(*composition, calculator, inputs, data, entry\_id=None, attribute=None, tag=None*)

Bases: MSONable

An lightweight Entry object containing key computed data for storing purpose.

**Parameters****composition**

[Composition] Composition of the entry. For flexibility, this can take the form of all the typical input taken by a Composition, including a {symbol: amt} dict, a string formula, and others.

**inputs**

[dict] An dict of parameters associated with the entry. Defaults to None.

**data**

[dict] An dict of any additional data associated with the entry. Defaults to None.

**entry\_id**

[obj] An optional id to uniquely identify the entry.

**attribute**

Optional attribute of the entry. This can be used to specify that the entry is a newly found compound, or to specify a particular label for the entry, or else ... Used for further analysis and plotting purposes. An attribute can be anything but must be MSONable.

**Attributes****number\_element**

## Methods

<code>as_dict()</code>	A JSON serializable dict representation of an object.
<code>from_dict(d)</code>	<p><b>param d</b> Dict representation.</p>
<code>to_json()</code>	Returns a json string representation of the MSONable object.
<code>unsafe_hash()</code>	Returns an hash of the current object.
<code>validate_monty_v1(_MSONable__input_value)</code>	Pydantic validator with correct signature for pydantic v1.x
<code>validate_monty_v2(_MSONable__input_value, _)</code>	Pydantic validator with correct signature for pydantic v2.x

### `as_dict()`

A JSON serializable dict representation of an object.

### classmethod `from_dict(d)`

#### Parameters

**d** – Dict representation.

#### Returns

MSONable class.

### property `number_element`

## `dpngen.database.run` module

`dpngen.database.run.db_run(args)`

`dpngen.database.run.parsing_gaussian(path, output='dpngen_db.json')`

`dpngen.database.run.parsing_pwscf(path, output='dpngen_db.json')`

`dpngen.database.run.parsing_vasp(path, config_info_dict, skip_init, output='dpngen_db.json', id_prefix=None)`

## `dpngen.database.vasp` module

`class dpngen.database.vasp.DPPotcar(symbols=None, functional='PBE', pp_file=None, pp_lists=None)`

Bases: MSONable



## Methods

<code>as_dict()</code>	A JSON serializable dict representation of an object.
<code>from_dict(d)</code>	<p><b>param d</b> Dict representation.</p>
<code>to_json()</code>	Returns a json string representation of the MSONable object.
<code>unsafe_hash()</code>	Returns an hash of the current object.
<code>validate_monty_v1(_MSONable__input_value)</code>	Pydantic validator with correct signature for pydantic v1.x
<code>validate_monty_v2(_MSONable__input_value, _)</code>	Pydantic validator with correct signature for pydantic v2.x

<b>from_file</b>
<b>write_file</b>

**as\_dict()**

A JSON serializable dict representation of an object.

**classmethod from\_dict(d)**

**Parameters**

**d** – Dict representation.

**Returns**

MSONable class.

**classmethod from\_file(filename)**

**write\_file(filename)**

**class** dpngen.database.vasp.**VaspInput**(*incar, poscar, potcar, kpoints=None, optional\_files=None, \*\*kwargs*)

Bases: dict, MSONable

Class to contain a set of vasp input objects corresponding to a run.

### Args:

*incar*: Incar object. *kpoints*: Kpoints object. *poscar*: Poscar object. *potcar*: Potcar object. *optional\_files*: Other input files supplied as a dict of {  
    *filename*: object}. The object should follow standard pymatgen conventions in implementing a `as_dict()` and `from_dict` method.

## Methods

<code>as_dict()</code>	A JSON serializable dict representation of an object.
<code>clear()</code>	
<code>copy()</code>	
<code>from_dict(d)</code>	<b>param d</b> Dict representation.
<code>from_directory(input_dir[, optional_files])</code>	Read in a set of VASP input from a directory.
<code>fromkeys(iterable[, value])</code>	Create a new dictionary with keys from iterable and values set to value.
<code>get(key[, default])</code>	Return the value for key if key is in the dictionary, else default.
<code>items()</code>	
<code>keys()</code>	
<code>pop(key[, default])</code>	If the key is not found, return the default if given; otherwise, raise a KeyError.
<code>popitem(/)</code>	Remove and return a (key, value) pair as a 2-tuple.
<code>setdefault(key[, default])</code>	Insert key with a value of default if key is not in the dictionary.
<code>to_json()</code>	Returns a json string representation of the MSONable object.
<code>unsafe_hash()</code>	Returns an hash of the current object.
<code>update([E, ]**F)</code>	If E is present and has a .keys() method, then does: for k in E: D[k] = E[k] If E is present and lacks a .keys() method, then does: for k, v in E: D[k] = v In either case, this is followed by: for k in F: D[k] = F[k]
<code>validate_monty_v1(_MSONable__input_value)</code>	Pydantic validator with correct signature for pydantic v1.x
<code>validate_monty_v2(_MSONable__input_value, _)</code>	Pydantic validator with correct signature for pydantic v2.x
<code>values()</code>	
<code>write_input([output_dir, ...])</code>	Write VASP input to a directory.

**as\_dict()**

A JSON serializable dict representation of an object.

**classmethod from\_dict(d)****Parameters**

**d** – Dict representation.

**Returns**

MSONable class.

**static from\_directory(input\_dir, optional\_files=None)**

Read in a set of VASP input from a directory. Note that only the standard INCAR, POSCAR, POTCAR and KPOINTS files are read unless optional\_filenames is specified.

**Parameters**

**input\_dir**

[str] Directory to read VASP input from.

**optional\_files**

[dict] Optional files to read in as well as a dict of {filename: Object type}. Object type must have a static method from\_file.

**write\_input**(*output\_dir='.', make\_dir\_if\_not\_present=True*)

Write VASP input to a directory.

**Parameters****output\_dir**

[str] Directory to write to. Defaults to current directory (“.”).

**make\_dir\_if\_not\_present**

[bool] Create the directory if not present. Defaults to True.

**dpngen.dispatcher package****Submodules****dpngen.dispatcher.Dispatcher module**

**dpngen.dispatcher.Dispatcher.make\_submission**(*mdata\_machine, mdata\_resources, commands, work\_path, run\_tasks, group\_size, forward\_common\_files, forward\_files, backward\_files, outlog, errlog*)

**dpngen.dispatcher.Dispatcher.make\_submission\_compat**(*machine: dict, resources: dict, commands: list[str], work\_path: str, run\_tasks: list[str], group\_size: int, forward\_common\_files: list[str], forward\_files: list[str], backward\_files: list[str], outlog: str = 'log', errlog: str = 'err', api\_version: str = '1.0'*) → None

Make submission with compatibility of both dispatcher API v0 and v1.

If *api\_version* is less than 1.0, raise RuntimeError. If *api\_version* is large than 1.0, use *make\_submission*.

**Parameters****machine**

[dict] machine dict

**resources**

[dict] resource dict

**commands**

[list[str]] list of commands

**work\_path**

[str] working directory

**run\_tasks**

[list[str]] list of paths to running tasks

**group\_size**

[int] group size

**forward\_common\_files**

[list[str]] forwarded common files shared for all tasks

**forward\_files**

[list[str]] forwarded files for each task

**backward\_files**

[list[str]] backwarded files for each task

**outlog**  
[str, default=log] path to log from stdout  
**errlog**  
[str, default=err] path to log from stderr  
**api\_version**  
[str, default=1.0] API version. 1.0 is required

`dpngen.dispatcher.Dispatcher.mdata_arginfo()` → `list[Argument]`

This method generates arginfo for a single mdata.

A submission requires the following keys: command, machine, and resources.

**Returns**  
`list[Argument]`  
arginfo

## **dpngen.generator package**

### **Subpackages**

#### **dpngen.generator.lib package**

### **Submodules**

#### **dpngen.generator.lib.abacus\_scf module**

`dpngen.generator.lib.abacus_scf.get_abacus_STRU(STRU, INPUT=None, n_ele=None)`  
`dpngen.generator.lib.abacus_scf.get_abacus_input_parameters(INPUT)`  
`dpngen.generator.lib.abacus_scf.get_additional_from_STRU(geometry_inlines, nele)`  
`dpngen.generator.lib.abacus_scf.get_mass_from_STRU(geometry_inlines, atom_names)`  
`dpngen.generator.lib.abacus_scf.get_natoms_from_stru(geometry_inlines)`  
`dpngen.generator.lib.abacus_scf.make_abacus_scf_input(fp_params, extra_file_path="")`  
`dpngen.generator.lib.abacus_scf.make_abacus_scf_kpt(fp_params)`  
`dpngen.generator.lib.abacus_scf.make_abacus_scf_stru(sys_data, fp_pp_files, fp_orb_files=None,  
fp_dpks_descriptor=None, fp_params=None,  
type_map=None, pporb="")`  
`dpngen.generator.lib.abacus_scf.make_kspacing_kpoints_stru(stru, kspacing)`  
`dpngen.generator.lib.abacus_scf.make_supercell_abacus(from_struct, super_cell)`

**dpgen.generator.lib.calypso\_check\_outcar module****dpgen.generator.lib.calypso\_run\_model\_devi module****dpgen.generator.lib.calypso\_run\_opt module****dpgen.generator.lib.cp2k module**

`dpgen.generator.lib.cp2k.iterdict(d, out_list, flag=None, indent=0)`

**Doc**

a recursive expansion of dictionary into cp2k input

**K**

current key

**V**

current value

**D**

current dictionary under expansion

**Flag**

used to record dictionary state. if flag is None,  
it means we are in top level dict. flag is a string. :indent: intent for current section.

`dpgen.generator.lib.cp2k.make_cp2k_input(sys_data, fp_params)`

`dpgen.generator.lib.cp2k.make_cp2k_input_from_external(sys_data, exinput_path)`

`dpgen.generator.lib.cp2k.make_cp2k_xyz(sys_data)`

`dpgen.generator.lib.cp2k.update_dict(old_d, update_d)`

A method to recursive update dict :old\_d: old dictionary :update\_d: some update value written in dictionary form.

**dpgen.generator.lib.cvasp module**

`dpgen.generator.lib.cvasp.runvasp(cmd, opt=False, max_errors=3, backup=False, auto_gamma=False, auto_npar=False, ediffg=-0.05)`

Cmd example: `cmd=['mpirun', '-np', '32', '-machinefile', 'hosts', 'vasp_std']`.

**dpgen.generator.lib.ele\_temp module**

`class dpgen.generator.lib.ele_temp.NBandsEsti(test_list)`

Bases: object

## Methods

<b>predict</b>
<b>save</b>

**predict**(*target\_dir*, *tolerance*=0.5)

**save**(*fname*)

### **dpgen.generator.lib.gaussian module**

dpgen.generator.lib.gaussian.**detect\_multiplicity**(*symbols*)

dpgen.generator.lib.gaussian.**make\_gaussian\_input**(*sys\_data*, *fp\_params*)

dpgen.generator.lib.gaussian.**take\_cluster**(*old\_conf\_name*, *type\_map*, *idx*, *jdata*)

### **dpgen.generator.lib.lammps module**

dpgen.generator.lib.lammps.**get\_all\_dumped\_forces**(*file\_name*)

dpgen.generator.lib.lammps.**get\_dumped\_forces**(*file\_name*)

dpgen.generator.lib.lammps.**make\_lammps\_input**(*ensemble*, *conf\_file*, *graphs*, *nsteps*, *dt*, *neidelay*, *trj\_freq*,  
*mass\_map*, *temp*, *jdata*, *tau\_t*=0.1, *pres*=None,  
*tau\_p*=0.5, *pka\_e*=None, *ele\_temp\_f*=None,  
*ele\_temp\_a*=None, *max\_seed*=1000000, *nopbc*=False,  
*deepmd\_version*='0.1', *nbeads*=None)

### **dpgen.generator.lib.make\_calypso module**

dpgen.generator.lib.make\_calypso.**make\_calypso\_input**(*nameofatoms*, *numberofatoms*, *numberofformula*,  
*volume*, *distanceofion*, *psoratio*, *popsze*,  
*maxstep*, *icode*, *split*, *vsc*, *maxnumatom*,  
*ctrlrange*, *pstress*, *fmax*)

dpgen.generator.lib.make\_calypso.**write\_model\_devi\_out**(*devi*, *fname*)

### **dpgen.generator.lib.parse\_calypso module**

### **dpgen.generator.lib.pwmat module**

dpgen.generator.lib.pwmat.**input\_upper**(*dinput*)

dpgen.generator.lib.pwmat.**make\_pwmat\_input\_dict**(*node1*, *node2*, *atom\_config*, *ecut*, *e\_error*, *rho\_error*,  
*icmix*=None, *smearing*=None, *sigma*=None,  
*kspacing*=0.5, *flag\_symm*=None)

`dpgen.generator.lib.pwmat.make_pwmat_input_user_dict(fp_params)`

`dpgen.generator.lib.pwmat.write_input_dict(input_dict)`

### **dpgen.generator.lib.pwscf module**

`dpgen.generator.lib.pwscf.cvt_1frame(fin, fout)`

`dpgen.generator.lib.pwscf.get_atom_types(lines)`

`dpgen.generator.lib.pwscf.get_block(lines, keyword, skip=0)`

`dpgen.generator.lib.pwscf.get_cell(lines)`

`dpgen.generator.lib.pwscf.get_coords(lines)`

`dpgen.generator.lib.pwscf.get_energy(lines)`

`dpgen.generator.lib.pwscf.get_force(lines)`

`dpgen.generator.lib.pwscf.get_natoms(lines)`

`dpgen.generator.lib.pwscf.get_stress(lines, cells)`

`dpgen.generator.lib.pwscf.get_types(lines)`

`dpgen.generator.lib.pwscf.make_pwscf_01_runctrl_dict(sys_data, idict)`

`dpgen.generator.lib.pwscf.make_pwscf_input(sys_data, fp_pp_files, fp_params, user_input=True)`

### **dpgen.generator.lib.run\_calypso module**

calypso as model devi engine: 1. gen\_structures 2. analysis 3. model devi.

`dpgen.generator.lib.run_calypso.analysis(iter_index, jdata, calypso_model_devi_path)`

`dpgen.generator.lib.run_calypso.gen_main(iter_index, jdata, mdata, caly_run_opt_list, gen_idx)`

`dpgen.generator.lib.run_calypso.gen_structures(iter_index, jdata, mdata, caly_run_path, current_idx, length_of_caly_runopt_list)`

`dpgen.generator.lib.run_calypso.run_calypso_model_devi(iter_index, jdata, mdata)`

### **dpgen.generator.lib.siesta module**

`dpgen.generator.lib.siesta.make_siesta_input(sys_data, fp_pp_files, fp_params)`

### dpgen.generator.lib.utils module

`dpgen.generator.lib.utils.cmd_append_log(cmd, log_file)`  
`dpgen.generator.lib.utils.copy_file_list(file_list, from_path, to_path)`  
`dpgen.generator.lib.utils.create_path(path)`  
`dpgen.generator.lib.utils.log_iter(task, ii, jj)`  
`dpgen.generator.lib.utils.log_task(message)`  
`dpgen.generator.lib.utils.make_iter_name(iter_index)`  
`dpgen.generator.lib.utils.record_iter(record, ii, jj)`  
`dpgen.generator.lib.utils.repeat_to_length(string_to_expand, length)`  
`dpgen.generator.lib.utils.replace(file_name, pattern, subst)`  
`dpgen.generator.lib.utils.symlink_user_forward_files(mdata, task_type, work_path, task_format=None)`  
Symlink user-defined forward\_common\_files Current path should be work\_path, such as 00.train.

#### Parameters

##### **mdata**

[dict] machine parameters

##### **task\_type**

[str] task\_type, such as “train”

##### **work\_path**

[str] work\_path, such as “iter.000001/00.train”

##### **task\_format**

[dict] formats of tasks

#### Returns

None

### dpgen.generator.lib.vasp module

`dpgen.generator.lib.vasp.incar_upper(dincar)`  
`dpgen.generator.lib.vasp.make_vasp_incar_user_dict(fp_params)`  
`dpgen.generator.lib.vasp.write_incar_dict(incar_dict)`

### Submodules

#### dpgen.generator.arginfo module

`dpgen.generator.arginfo.basic_args()` → list[Argument]  
`dpgen.generator.arginfo.data_args()` → list[Argument]  
`dpgen.generator.arginfo.fp_args()` → list[Argument]



`dpgen.generator.arginfo.fp_style_abacus_args()` → list[Argument]

`dpgen.generator.arginfo.fp_style_amber_diff_args()` → list[Argument]

Arguments for FP style amber/diff.

**Returns**

**list[dargs.Argument]**

list of amber/diff fp style arguments

`dpgen.generator.arginfo.fp_style_cp2k_args()` → list[Argument]

`dpgen.generator.arginfo.fp_style_custom_args()` → list[Argument]

Arguments for FP style custom.

**Returns**

**list[dargs.Argument]**

list of custom fp style arguments

`dpgen.generator.arginfo.fp_style_gaussian_args()` → list[Argument]

Gaussian fp style arguments.

**Returns**

**list[dargs.Argument]**

list of Gaussian fp style arguments

`dpgen.generator.arginfo.fp_style_pwscf_args()` → list[Argument]

Arguments for FP style pwscf (Quantum Espresso).

**Returns**

**list[dargs.Argument]**

list of pwscf fp style arguments

`dpgen.generator.arginfo.fp_style_siesta_args()` → list[Argument]

`dpgen.generator.arginfo.fp_style_variant_type_args()` → Variant

`dpgen.generator.arginfo.fp_style_vasp_args()` → list[Argument]

`dpgen.generator.arginfo.model_devi_amber_args()` → list[Argument]

Amber engine arguments.

`dpgen.generator.arginfo.model_devi_args()` → list[Variant]

`dpgen.generator.arginfo.model_devi_jobs_args()` → list[Argument]

`dpgen.generator.arginfo.model_devi_jobs_rev_mat_args()` → Argument

`dpgen.generator.arginfo.model_devi_jobs_template_args()` → Argument

`dpgen.generator.arginfo.model_devi_lmp_args()` → list[Argument]

`dpgen.generator.arginfo.run_jdata_arginfo()` → Argument

Argument information for dpngen run mdata.

**Returns**

**Argument**

argument information

`dpgen.generator.arginfo.run_mdata_arginfo()` → Argument

Generate arginfo for dpngen run mdata.

**Returns**

**Argument**

arginfo

`dpgen.generator.arginfo.training_args()` → list[Argument]

Traning arguments.

**Returns**

**list[dargs.Argument]**

List of training arguments.

## **dpgen.generator.run module**

init: data iter:

00.train 01.model\_devi 02.vasp 03.data.

`dpgen.generator.run.check_bad_box(conf_name, criteria, fmt='lammps/dump')`

`dpgen.generator.run.check_cluster(conf_name, fp_cluster_vacuum, fmt='lammps/dump')`

`dpgen.generator.run.copy_model(numb_model, prv_iter_index, cur_iter_index)`

`dpgen.generator.run.detect_batch_size(batch_size, system=None)`

`dpgen.generator.run.dump_to_deepmd_raw(dump, deepmd_raw, type_map, fmt='gromacs/gro',  
charge=None)`

`dpgen.generator.run.expand_idx(in_list)`

`dpgen.generator.run.expand_matrix_values(target_list, cur_idx=0)`

`dpgen.generator.run.find_only_one_key(lmp_lines, key)`

`dpgen.generator.run.gen_run(args)`

`dpgen.generator.run.get_atomic_masses(atom)`

`dpgen.generator.run.get_job_names(jdata)`

`dpgen.generator.run.get_nframes(system)`

`dpgen.generator.run.get_sys_index(task)`

`dpgen.generator.run.make_fp(iter_index, jdata, mdata)`

Select the candidate strutures and make the input file of FP calculation.

**Parameters**

**iter\_index**

[int] iter index

**jdata**

[dict] Run parameters.

**mdata**

[dict] Machine parameters.

`dpgen.generator.run.make_fp_abacus_scf(iter_index, jdata)`

`dpgen.generator.run.make_fp_amber_diff(iter_index: int, jdata: dict)`

Run amber twice to calculate high-level and low-level potential, and then generate difference between them.

Besides AMBER, one needs to install *dpamber* package, which is avaiable at <https://github.com/njzjz/dpamber>

Currently, it should be used with the AMBER model\_devi driver.

**Parameters**

**iter\_index**

[int] iter index

**jdata**

[dict]

**Run parameters. The following parameters are used in this method:****mdin\_prefix**

[str] The path prefix to AMBER mdin files

**qm\_region**

[list[str]] AMBER mask of the QM region. Each mask maps to a system.

**qm\_charge**

[list[int]] Charge of the QM region. Each charge maps to a system.

**high\_level**

[str] high level method

**low\_level**

[str] low level method

**fp\_params**

[dict]

**This parameters includes:****high\_level\_mdin**

[str] High-level AMBER mdin file. %qm\_theory%, %qm\_region%, and %qm\_charge% will be replace.

**low\_level\_mdin**

[str] Low-level AMBER mdin file. %qm\_theory%, %qm\_region%, and %qm\_charge% will be replace.

**parm7\_prefix**

[str] The path prefix to AMBER PARM7 files

**parm7**

[list[str]] List of paths to AMBER PARM7 files. Each file maps to a system.

## References

[1]

`dpngen.generator.run.make_fp_calculation(iter_index, jdata, mdata)`

Make the input file of FP calculation.

**Parameters****iter\_index**

[int] iter index

**jdata**

[dict] Run parameters.

**mdata**

[dict] Machine parameters.

`dpngen.generator.run.make_fp_cp2k(iter_index, jdata)``dpngen.generator.run.make_fp_custom(iter_index, jdata)`

Make input file for customized FP style.

Convert the POSCAR file to custom format.

**Parameters****iter\_index**

[int] iter index

**jdata**

[dict] Run parameters.

```
dpgen.generator.run.make_fp_gaussian(iter_index, jdata)
dpgen.generator.run.make_fp_pwmat(iter_index, jdata)
dpgen.generator.run.make_fp_pwscf(iter_index, jdata)
dpgen.generator.run.make_fp_siesta(iter_index, jdata)
dpgen.generator.run.make_fp_task_name(sys_idx, counter)
dpgen.generator.run.make_fp_vasp(iter_index, jdata)
dpgen.generator.run.make_fp_vasp_cp_cvasp(iter_index, jdata)
dpgen.generator.run.make_fp_vasp_incar(iter_index, jdata, nbands_esti=None)
dpgen.generator.run.make_fp_vasp_kp(iter_index, jdata)
dpgen.generator.run.make_model_devi(iter_index, jdata, mdata)
dpgen.generator.run.make_model_devi_conf_name(sys_idx, conf_idx)
dpgen.generator.run.make_model_devi_task_name(sys_idx, task_idx)
dpgen.generator.run.make_pwmat_input(jdata, filename)
dpgen.generator.run.make_train(iter_index, jdata, mdata)
dpgen.generator.run.make_vasp_incar(jdata, filename)
dpgen.generator.run.make_vasp_incar_ele_temp(jdata, filename, ele_temp, nbands_esti=None)
dpgen.generator.run.parse_cur_job(cur_job)
dpgen.generator.run.parse_cur_job_revmat(cur_job, use_plm=False)
dpgen.generator.run.parse_cur_job_sys_revmat(cur_job, sys_idx, use_plm=False)
dpgen.generator.run.poscar_natoms(lines)
dpgen.generator.run.poscar_shuffle(poscar_in, poscar_out)
dpgen.generator.run.poscar_to_conf(poscar, conf)
dpgen.generator.run.post_fp(iter_index, jdata)
dpgen.generator.run.post_fp_abacus_scf(iter_index, jdata)
dpgen.generator.run.post_fp_amber_diff(iter_index, jdata)
dpgen.generator.run.post_fp_check_fail(iter_index, jdata, rfailed=None)
dpgen.generator.run.post_fp_cp2k(iter_index, jdata, rfailed=None)
```

`dpgen.generator.run.post_fp_custom(iter_index, jdata)`

Post fp for custom fp. Collect data from user-defined *output\_fn*.

**Parameters**

**iter\_index**

[int] The index of the current iteration.

**jdata**

[dict] The parameter data.

`dpgen.generator.run.post_fp_gaussian(iter_index, jdata)`

`dpgen.generator.run.post_fp_pwmat(iter_index, jdata, rfailed=None)`

`dpgen.generator.run.post_fp_pwscf(iter_index, jdata)`

`dpgen.generator.run.post_fp_siesta(iter_index, jdata)`

`dpgen.generator.run.post_fp_vasp(iter_index, jdata, rfailed=None)`

`dpgen.generator.run.post_model_devi(iter_index, jdata, mdata)`

`dpgen.generator.run.post_train(iter_index, jdata, mdata)`

`dpgen.generator.run.revise_by_keys(lmp_lines, keys, values)`

`dpgen.generator.run.revise_lmp_input_dump(lmp_lines, trj_freq, model_devi_merge_traj=False)`

`dpgen.generator.run.revise_lmp_input_model(lmp_lines, task_model_list, trj_freq, deepmd_version='1')`

`dpgen.generator.run.revise_lmp_input_plm(lmp_lines, in_plm, out_plm='output.plumed')`

`dpgen.generator.run.run_fp(iter_index, jdata, mdata)`

`dpgen.generator.run.run_fp_inner(iter_index, jdata, mdata, forward_files, backward_files, check_fn, log_file='fp.log', forward_common_files=[])`

`dpgen.generator.run.run_iter(param_file, machine_file)`

`dpgen.generator.run.run_md_model_devi(iter_index, jdata, mdata)`

`dpgen.generator.run.run_model_devi(iter_index, jdata, mdata)`

`dpgen.generator.run.run_train(iter_index, jdata, mdata)`

`dpgen.generator.run.set_version(mdata)`

`dpgen.generator.run.sys_link_fp_vasp_pp(iter_index, jdata)`

`dpgen.generator.run.update_mass_map(jdata)`

## dpgen.remote package

### Submodules

#### dpgen.remote.decide\_machine module

`dpgen.remote.decide_machine.convert_mdata(mdata, task_types=['train', 'model_devi', 'fp'])`

Convert mdata for DP-GEN main process. New conversion is like `mdata["fp"]["machine"]`, DP-GEN needs `mdata["fp_machine"]`.

Notice that we deprecate the function which can automatically select one most available machine, since this function was only used by Angus, and only supports for Slurm. In the future this can be implemented.

**Parameters**

**mdata**

[dict] Machine parameters to be converted.

**task\_types**

[list of string] Type of tasks, default is ["train", "model\_devi", "fp"]

**Returns**

**dict**

mdata converted

## dpgen.simplify package

### Submodules

#### dpgen.simplify.arginfo module

`dpgen.simplify.arginfo.fp_args()` → list[Argument]

Generate arginfo for fp.

**Returns**

**List[Argument]**

arginfo

`dpgen.simplify.arginfo.fp_style_variant_type_args()` → Variant

Generate variant for fp style variant type.

**Returns**

**Variant**

variant for fp style

`dpgen.simplify.arginfo.general_simplify_arginfo()` → Argument

General simplify arginfo.

**Returns**

**Argument**

arginfo

`dpgen.simplify.arginfo.simplify_jdata_arginfo()` → Argument

Generate arginfo for dpgen simplify jdata.

**Returns**

**Argument**

arginfo

`dpgen.simplify.arginfo.simplify_mdata_arginfo()` → Argument

Generate arginfo for dpgen simplify mdata.

**Returns**

**Argument**

arginfo

## dpngen.simplify.simplify module

Simplify dataset (minimize the dataset size).

Init: pick up init data from dataset randomly

Iter: 00: train models (same as generator) 01: calculate model deviations of the rest dataset, pick up data with proper model deviation 02: fp (optional, if the original dataset do not have fp data, same as generator)

`dpngen.simplify.simplify.gen_simplify(args)`

`dpngen.simplify.simplify.get_multi_system(path: str | list[str], jdata: dict) → MultiSystems`

Get MultiSystems from a path or list of paths.

Both NumPy and HDF5 formats are supported. For details of two formats, refer to DeePMD-kit documentation.

If *labeled* in *jdata* is True, returns MultiSystems with LabeledSystem. Otherwise, returns MultiSystems with System.

### Parameters

#### path

[str or list of str] path or list of paths to the dataset

#### jdata

[dict] parameters which may contain *labeled* key

### Returns

#### dpdata.MultiSystems

MultiSystems with LabeledSystem or System

`dpngen.simplify.simplify.get_system_cls(jdata)`

`dpngen.simplify.simplify.init_model(iter_index, jdata, mdata)`

`dpngen.simplify.simplify.init_pick(iter_index, jdata, mdata)`

Pick up init data from dataset randomly.

`dpngen.simplify.simplify.make_fp(iter_index, jdata, mdata)`

`dpngen.simplify.simplify.make_fp_configs(iter_index, jdata)`

`dpngen.simplify.simplify.make_fp_labeled(iter_index, jdata)`

`dpngen.simplify.simplify.make_model_devi(iter_index, jdata, mdata)`

Calculate the model deviation of the rest idx.

`dpngen.simplify.simplify.post_model_devi(iter_index, jdata, mdata)`

Calculate the model deviation.

`dpngen.simplify.simplify.run_iter(param_file, machine_file)`

Init (iter 0): init\_pick.

tasks (iter > 0): 00 make\_train (same as generator) 01 run\_train (same as generator) 02 post\_train (same as generator) 03 make\_model\_devi 04 run\_model\_devi 05 post\_model\_devi 06 make\_fp 07 run\_fp (same as generator) 08 post\_fp (same as generator)

`dpngen.simplify.simplify.run_model_devi(iter_index, jdata, mdata)`

Submit dp test tasks.

## dpngen.tools package

### Submodules

#### dpngen.tools.auto\_gen\_param module

```
class dpngen.tools.auto_gen_param.Iteration(temps, nsteps_list=[500, 500, 1000, 1000, 3000, 3000, 6000, 6000], sub_iteration_num=8, ensemble='npt', press=[1.0, 10.0, 100.0, 1000.0, 5000.0, 10000.0, 20000.0, 50000.0], trj_freq=10)
```

Bases: object

#### Attributes

`index_iteration`

#### Methods

<code>gen_sub_iter</code>
<code>register_iteration</code>
<code>register_sub_iteartion</code>

```
current_num_of_itearation = 0
```

```
current_num_of_sub_itearation = 0
```

```
gen_sub_iter(system_list)
```

```
property index_iteration
```

```
classmethod register_iteration()
```

```
classmethod register_sub_iteartion()
```

```
class dpngen.tools.auto_gen_param.System(system_prefix="")
```

Bases: object

#### Attributes

`index_system`

#### Methods

<code>add_sub_system</code>
<code>get_sub_system</code>
<code>register_sub_system</code>
<code>register_system</code>

```
add_sub_system(idx2, files_list)
```

```
current_num_of_sub_systems = 0
```

```
current_num_of_system = 0
```



```

    get_sub_system()

    property index_system

    classmethod register_sub_system()

    classmethod register_system()

dpgen.tools.auto_gen_param.auto_gen_param(args)

dpgen.tools.auto_gen_param.default_map_generator(map_list=[1, 1, 2, 2, 2, 4, 4, 4], data_list=None)

dpgen.tools.auto_gen_param.default_temps_generator(melt_point, temps_interval=0.1, num_temps=5)

dpgen.tools.auto_gen_param.get_basic_param_json(melt_point, out_param_filename='param_basic.json',
                                                  scan_dir='.', file_name='POSCAR',
                                                  init_file_name='type.raw', min_allow_files_num=16,
                                                  map_list=[1, 1, 2, 2, 2, 4, 4, 4], meta_iter_num=4,
                                                  sub_iteration_num=8, map_iterator=None,
                                                  nsteps_list=[500, 500, 1000, 1000, 3000, 3000, 6000,
                                                                6000], press=[1.0, 10.0, 100.0, 1000.0, 5000.0,
                                                                10000.0, 20000.0, 50000.0], temps_iterator=None,
                                                  ensemble='npt', trj_freq=10, temps_interval=0.1,
                                                  num_temps=5)

dpgen.tools.auto_gen_param.get_init_data_sys(scan_dir='.', init_file_name='type.raw')

dpgen.tools.auto_gen_param.get_model_devi_jobs(melt_point, system_list, nsteps_list=[500, 500, 1000,
                                                                                   1000, 3000, 3000, 6000, 6000], press=[1.0, 10.0,
                                                                                   100.0, 1000.0, 5000.0, 10000.0, 20000.0, 50000.0],
                                                                                   meta_iter_num=4, sub_iteration_num=8,
                                                                                   temps_iterator=None, ensemble='npt', trj_freq=10,
                                                                                   temps_interval=0.1, num_temps=5)

dpgen.tools.auto_gen_param.get_sys_configs(system_list)

dpgen.tools.auto_gen_param.get_system_list(system_dict, map_list=[1, 1, 2, 2, 2, 4, 4, 4],
                                           meta_iter_num=4, sub_iteration_num=8,
                                           map_iterator=None, file_name='POSCAR')

:Exmaple [['000000', '000001'], ['00000[2-9]'], ['00001?', '000020'],]

dpgen.tools.auto_gen_param.scan_files(scan_dir='.', file_name='POSCAR', min_allow_files_num=20)

```

### dpgen.tools.collect\_data module

```

dpgen.tools.collect_data.collect_data(target_folder, param_file, output, verbose=True)

dpgen.tools.collect_data.file_len(fname)

```

### dpgen.tools.relabel module

```
dpgen.tools.relabel.copy_pp_files(tdir, fp_pp_path, fp_pp_files)
dpgen.tools.relabel.create_init_tasks(target_folder, param_file, output, fp_json, verbose=True)
dpgen.tools.relabel.create_tasks(target_folder, param_file, output, fp_json, verbose=True, numb_iter=-1)
dpgen.tools.relabel.get_lmp_info(input_file)
dpgen.tools.relabel.link_pp_files(tdir, fp_pp_path, fp_pp_files)
dpgen.tools.relabel.make_pwscf(tdir, fp_params, mass_map, fp_pp_path, fp_pp_files, user_input)
dpgen.tools.relabel.make_siesta(tdir, fp_params, fp_pp_path, fp_pp_files)
dpgen.tools.relabel.make_vasp(tdir, fp_params)
dpgen.tools.relabel.make_vasp_incar(tdir, fp_incar)
```

### dpgen.tools.run\_report module

```
dpgen.tools.run_report.run_report(args)
```

### dpgen.tools.stat\_iter module

```
dpgen.tools.stat_iter.stat_iter(target_folder, param_file='param.json', verbose=True, mute=False)
```

### dpgen.tools.stat\_sys module

```
dpgen.tools.stat_sys.ascii_hist(count)
dpgen.tools.stat_sys.run_report(args)
dpgen.tools.stat_sys.stat_sys(target_folder, param_file='param.json', verbose=True, mute=False)
```

### dpgen.tools.stat\_time module

```
dpgen.tools.stat_time.stat_time(target_folder, param_file='param.json', verbose=True, mute=False)
```

## 10.1.2 Submodules

### 10.1.3 dpgen.arginfo module

```
dpgen.arginfo.general_mdata_arginfo(name: str, tasks: tuple[str]) → Argument
```

Generate arginfo for general mdata.

#### Parameters

**name**

[str] mdata name

**tasks**  
 [tuple[str]] tuple of task keys, e.g. (“train”, “model\_devi”, “fp”)  
**Returns**  
**Argument**  
 arginfo

### 10.1.4 dpngen.gui module

DP-GUI entrypoint.

`dpngen.gui.start_dpgui(args: Namespace)`

Host DP-GUI server.

**Parameters**

**args**

[argparse.Namespace] Arguments from argparse.

**Raises**

**ModuleNotFoundError**

The dpgui package is not installed

### 10.1.5 dpngen.main module

`dpngen.main.main()`

`dpngen.main.main_parser()` → ArgumentParser

Returns parser for *dpngen* command.

**Returns**

**argparse.ArgumentParser**

parser for *dpngen* command

### 10.1.6 dpngen.util module

`dpngen.util.box_center(ch='', fill=' ', sp='|')`

Put the string at the center of | |.

`dpngen.util.convert_training_data_to_hdf5(input_files: list[str], h5_file: str)`

Convert training data to HDF5 format and update the input files.

**Parameters**

**input\_files**

[list of str] DeePMD-kit input file names

**h5\_file**

[str] HDF5 file name

`dpngen.util.expand_sys_str(root_dir: str | Path) → list[str]`

Recursively iterate over directories taking those that contain *type.raw* file.

If root\_dir is a file but not a directory, it will be assumed as an HDF5 file.

**Parameters**

**root\_dir**

[Union[str, Path]] starting directory

**Returns**

**List[str]**

list of string pointing to system directories

**Raises****RuntimeError**

No system was found in the directory

`dpgen.util.load_file(filename: str | PathLike) → dict`

Load data from a JSON or YAML file.

**Parameters****filename**

[str or os.PathLike] The filename to load data from, whose suffix should be .json, .yaml, or .yaml

**Returns****dict**

The data loaded from the file

**Raises****ValueError**

If the file format is not supported

`dpgen.util.normalize(arginfo: Argument, data: dict, strict_check: bool = True) → dict`

Normalize and check input data.

**Parameters****arginfo**

[dargs.Argument] argument information

**data**

[dict] input data

**strict\_check**

[bool, default=True] strict check data or not

**Returns****dict**

normalized data

`dpgen.util.sepline(ch='-', sp='-', screen=False)`

Seperate the output by ‘-‘.

`dpgen.util.set_directory(path: Path)`

Sets the current working path within the context.

**Parameters****path**

[Path] The path to the cwd

**Yields****None**

## Examples

```
>>> with set_directory("some_path"):  
...     do_something()
```

`dpgen.util.setup_ele_temp(atomic: bool)`

Set electronic temperature as required input data.

**Parameters****atomic**

[bool] Whether to use atomic temperature or frame temperature

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- genindex
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