

ELASTOOL Users Manual

AN AUTOMATED TOOLKIT FOR ELASTIC CONSTANTS CALCULATION

VERSION: 1.0.2

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1 Introduction

1.1 About ELASTOOL

ELASTOOL is an automated toolkit for calculating the second-order elastic constants (SOECs) of any crystal systems belonging to two- and three-dimensional. It can utilize three kinds of strain-matrix sets, the high efficiency strain-matrix sets (OHESS) [?], the universal linear-independent coupling strains (ULICS) [?] and the all single-element strain-matrix sets (ASESS) [?] to calculate the SOECs automatically. In an automatic manner, ELASTOOL can deal with both zero- and high-temperature elastic constants.

Presently, ELASTOOL interfaces to VASP package for calculating the accurate stresses of strained crystal. But the interfaces to other DFT packages can also be easily implemented.

1.2 Features

ELASTOOL has many features including

- Very easy to use (installation and run);
- High efficiency;
- Automated flow of the SOECs calculation;
- Three kinds of strain-matrix sets, the OHESS, the ASESS, and the ULICS;
- Zero-temperature SOECs;
- High-temperature SOECs;
- Interfaced to VASP.

2 Installation and run

ELASTOOL is based on Python and its installation is very easy. But before the installation of ELASTOOL, the necessary libraries should be installed first. The following packages are required:

2.1 Installation requirements

ELASTOOL depends on Python 3. The following packages are required:

Python3.5 or later.

- NumPy.
- Spglib.
- ASE.
- Pandas.
- VASP.

As for Spglib, please install spglib for the structruces' space group determination.

2.2 Installation

In the Python 3 enviroment, the necessary libraries can be installed via pip command, e.g., `pip install numpy`. For the construction of the environment and libraries, the `miniconda3` management platform of Python packages is highly recommended. After installing `miniconda3`, the basic Python 3 language environment is cocnstructed and the other libraries can be installed via either `conda` or `pip` commands. For example, one can install `numpy` via `conda install -c conda-forge numpy`.

2.3 Run

To run ELASTOOL, one only needs to execute `elastool` in the working directory. ELASTOOL will automatically prepare necessary files for calculating the stresses of crystal under deformation, and then call VASP to optimize initial crystal structure and calculate the stresses for each deformaton defined by `OHESS`, `ASESS`, or `ULICS`. Finally, `ElasTool` analyzes stress-strain relationship according Hooke's law and calculate all the elastic constants.

2.3.1 Interface to VASP

To run ELASTOOL with VASP, one need provide `elastic.in`, `INCARs`, and `POTCAR-XX` files for each kind of atom, where `XX` is the short name of each atom. For example, you can use `POTCAR-Mg`, and `POTCAR-O` for `MgO`.

2.4 Input files

ELASTOOL needs one main input file for setting the calculation details of elastic constants. The main input file is named `elatoool.in`. The crystal structure file is provided either in `POSCAR` or `cif` format for reading in the structure information of the crystal. For VASP stress tensor calculations, `INCARs`, `KPOINTS-static`, `KPOINTS-dynamic`, and `POTCAR-XX` files are also necessary. In the `INCARs` file, several `INCAR` files of VASP are collected for optimization, the static calculation, and the molecular dynamics simulations for high-temperature elastic constants. The `KPOINTS-static` file is for the structure optimization and the static calculation of stress tensors. The `KPOINTS-dynamic` file is for the calculation of high-temperature elastic constants using molecular dynamics. `XX` in the `POTCAR` file name is the abbreviated name for an element of the crystal.

3 Input parameters

There are totally 10 controlling parameters of ELASTOOL, as listed in Table.1. The `run_mode` sets the running mode for ElastoTool, 1 for automatic run, 2 for pre-processing, and 3 for post-processing. If `run_mode` = 2 or 3, one should ensure the structure has already been optimized at fixed pressure or volume, *i.e.* both the CONTCAR and OUTCAR files are in `./OPT` directory. In running mode 2, ELASTOOL will directly prepare all the necessary files for calculating stress tensors. After all the stress tensors calculations are finished, run mode 3 can analyze the output files and extract stress tensors, and then fit the first-order function to the stress-tensor data to obtain elastic constants. The `dimensional` defines the dimensional of the system, 2D or 3D. If the system is 2D, ELASTOOL supposes the layered sheet in the *xy*-plane. The `structure_file` specifies the original crystal structure file in POSCAR (.vasp) or cif (.cif) format. The `if_conventional_cell` determines the usage of primitive cell (no) or conventional cell (yes). The `method_stress_statistics` chooses the elastic constants calculation method, static or dynamic, static for 0 K elastic constants, dynamic for high-temperature. The static method uses the static stress to compute elastic constants, while the dynamic method deduces elastic constants from the thermal stresses obtained by molecular dynamics simulations. The `strains_matrix` defines the type of strain-matrix set, OHES, ASESS, or ULICS. The `strains_list` gives one or more strains for calculating stresses via the strain-matrix set of OHES, ASESS, or ULICS. The `repeat_num` controls how to build a supercell from the primitive or conventional cell defined by `if_conventional_cell` for the dynamic method. The `num_last_samples` is the number of last MD steps to average thermal stresses. The `parallel_submit_command` is the parallel submitting command of *ab initio* code, e.g. VASP.

Table 1. The controlling parameters and possible values of ELASTOOL

Parameters	Values
<code>run_mode</code>	1/2/3
<code>dimensional</code>	2D/3D
<code>structure_file</code>	file name ended with .vasp or .cif
<code>if_conventional_cell</code>	yes/no
<code>method_stress_statistics</code>	static/dynamic
<code>strains_matrix</code>	ohess/assess/ulics
<code>strains_list</code>	one or more numbers
<code>repeat_num</code>	3 integers
<code>num_last_samples</code>	1 integer
<code>parallel_submit_command</code>	DFT parallel run command

4 Example of run

The best way to learn ELASTOOL is to start from the examples. ELASTOOL can calculate zero-temperature and high-temperature elastic constants. The zero-temperature calculations can be conducted by the static stress computation. The high-temperature elastic constants can be derived by molecular dynamics simulations.

4.1 Zero-temperature elastic constants

We take the 0 K elastic constants calculation of diamond as the static example. The content of the input file elastool.in is as follows.

```
run_mode = 1
dimensional = 3D
structure_file = diamond.cif
if_conventional_cell = no
method_stress_statistics = static
strains_matrix = ohess
strains_list = -0.06 -0.03 0.03 0.06
#repeat_num = 1 1 1
#num_last_samples = 1
parallel_submit_command = mpirun -np 28 vasp544
```

4.2 High-temperature elastic constants

4.2.1 3D case

The high temperature elastic constants calculation of metal copper is the high-temperature example. We build a $3 \times 3 \times 3$ supercell from the conventional cell of face-centered-cubic of Cu and then perform long-time MD simulations defined in the INCAR-dynamic file. Because MD is very time consuming, there is only one strain of -0.06 is used. The last 500 MD steps are used to average thermal stresses.

```
run_mode = 1
dimensional = 3D
structure_file = CONTCAR.vasp
if_conventional_cell = yes
method_stress_statistics = dynamic
strains_matrix = ohess
strains_list = -0.06
repeat_num = 3 3 3
num_last_samples = 500
parallel_submit_command = mpirun -np 28 vasp544
```

4.2.2 2D case

```
run_mode = 1
```

```

dimensional = 2D
structure_file = CONTCAR.vasp
if_conventional_cell = no
method_stress_statistics = dynamic
strains_matrix = o Hess
strains_list = -0.06
repeat_num = 4 4 1
num_last_samples = 500
parallel_submit_command = mpirun -np 28 vasp544

```

4.3 Output files

The elastool.out file is the unique output file of ELASTOOL. It includes the calculated elastic constants data, the elastic moduli, the sound velocity, the Debye temperature, the elastic anisotropy, and the stability analysis results of the crystal structure based on Born elastic criteria. The printed information on screen of the diamond example is as follows.

Reading controlling parameters from elastool.in...

Calculating stresses using the OHESS strain matrices...

strain = -0.060

strain = -0.030

strain = 0.030

strain = 0.060

Fitting the first-order function to the collected

stress-strain data according to Hooke's law...

The final results are as follows:

```

+=====+
|This is a 3D Cubic lattice.      |
|_____|
|Mean Pressure = -0.08 GPa       |
|_____|
|Elastic constants:              |
|C11 = 1055.04 GPa               |
|C12 = 136.56 GPa                |
|C44 = 567.76 GPa                |
|_____|

```

```

|Elastic moduli:                                     |
|B_V = 442.72 GPa                                   |
|B_R = 442.72 GPa                                   |
|G_V = 524.36 GPa                                   |
|G_R = 518.73 GPa                                   |
|B_VRH = 442.72 GPa                                 |
|G_VRH = 521.54 GPa                                 |
|Young's modulus (E) = 1123.47 GPa                  |
|Possion's ratio (V) = 0.0771                      |
|-----|
|Sound velocity:                                     |
|V_S = 12.20 Km/s                                   |
|V_B = 11.24 Km/s                                   |
|V_P = 18.03 Km/s                                   |
|V_M = 13.31 Km/s                                   |
|-----|
|Debye temperature:                                 |
|T_D = 1761.03 K                                    |
|-----|
|Elastic anisotropy:                                |
|A_U = 0.0542                                        |
|A_C = 0.0054                                        |
|-----|
|Structure stability analysis...                     |
|This structure is mechanically STABLE.             |
+=====+
Results are also saved in the elastool.out file.
Well done! GOOD LUCK!

```

5 How to cite

Please cite the following article when you use ELASTOOL:

Z. L. Liu. ELASTOOL: An automated toolkit for elastic constants calculation (arxiv:2002.06535). 2020.

Z. L. Liu. High-efficiency calculation of elastic constants enhanced by the optimized strain-matrix sets (arxiv:2002.00005). 2020.

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Bibliography