

# **CGFD3D-elastic**

## **User Manual**

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# Contents

<b>Contents</b>	<b>1</b>
<b>1 Grid</b>	<b>2</b>
1.1 Layer Interp . . . . .	3
1.1.1 File Format (.gdlay) . . . . .	3
1.1.2 Example . . . . .	3
<b>2 How to give media parameters and run wave simulations for different media types</b>	<b>5</b>
2.1 Format of .md3grd . . . . .	6
2.2 Format of .md3lay . . . . .	7
2.3 Note about equivalent medium parameterization methods . . . . .	8
<b>3 Source</b>	<b>9</b>
3.1 JSON read source format . . . . .	10
3.2 analytic source format . . . . .	10
3.3 value source format . . . . .	11
<b>Copyright</b>	<b>13</b>
<b>Bibliography</b>	<b>14</b>

# Chapter 1

## Grid

run\_test.sh

---

```
"#grid_generation_method" : {  
  "#import" : "$GRID_DIR",  
  "#cartesian" : {  
    "origin" : [0.0, 0.0, -5900.0 ],  
    "interval" : [ 100.0, 100.0, 100.0 ]  
  },  
  "layer_interp" : {  
    "in_grid_layer_file" : "$EXEC_DIR/test/test_grid.gdlay",  
    "refine_factor" : [ 1, 1, 1 ],  
    "horizontal_start_index" : [ 3, 3 ],  
    "vertical_ToFreeSurf_resample_index" : 0  
  }  
},  
"is_export_grid" : 1,  
"grid_export_dir" : "$GRID_DIR",
```

---

In the configuration of grid, we can choose different grid generation methods.

(1) if `import` is selected, we can directly import the grid file from the outside. `$GRID_DIR` is the file path and name.

(2) if `cartesian` is selected, we can get the grid through interpolation with several given control interfaces. we can generate grids in Cartesian coordinates. Two parameters are required here, `origin` is the starting coordinate of the grid and `interval` is the grid spacing.

(3) if `layer_interp` is selected, we can get the grid through interpolation with several given control interfaces. there need to be a `.gdlay` file, and the file format is shown in Section 1.1.

`refine_factor` is resampling control parameters for interface grids. This parameter must be an integer and cannot be zero. eg: If the parameter is 1, it means no resampling; If this parameter is `-2` or `-3`, it means two or three times downsampling; If this parameter is 2 or 3, it means two or three times interpolating. Its value is greater than or equal to the number of ghost points

`horizontal_start_index` and `vertical_ToFreeSurf_resample_index` intercept sub-models from the interface control model.

`horizontal_start_index` is x and y start index of interface before resampling. Note: ensure that after resampling, three points must be set aside as ghost points outside the sub-model. eg: If you do not resample, the minimum value of this parameter is 3; If you do two or three times downsampling, the minimum value of this parameter is 6 or 10; If you do two or three times interpolating, the minimum value of this parameter is 2 or 1;

`vertical_ToFreeSurf_resample_index` is index to free surface. If this parameter is 0, it means selecting the submodel that contains the free surface; If this parameter is 40, it means selecting the submodel that intercepted at 40 nodes from the free surface in the resampled interface control model.

## 1.1 Layer Interp

The **layer\_interp** function interpolates the spatial grid of the model through the given interface.

### 1.1.1 File Format (.gdlay)

The following description ignores comment lines and blank lines.

- The first line is the number of interface.
- The second line is the cells of each layer.
- The third line is isequal of each layer.
- The fourth line is number of interface grids in X and Y directions.
- Each subsequent line is the X, Y and Z coordinates of the grids.

### 1.1.2 Example

A model with a horizontal interface can be given as:

test.gdlay			
4			
20	10	20	
1	0	1	
100	100		
100.0	100.0	-5000.0	
200.0	100.0	-5000.0	
300.0	100.0	-5000.0	
400.0	100.0	-5000.0	
...	...	...	

The given interface model have 4 interfaces.

There are 20 cells between the first interface and the second interface; there are 10 cells between the second interface and the third interface; there are 20 cells between the third interface and the fourth interface. The first interface is the deepest underground interface, and the last is the free surface interface.

The cell spacing between the first interface and the second interface are equal; the cell spacing between the third interface and the fourth interface are equal; the cell spacing between the second interface and the third interface is gradual. The number of interface grids in X is 100 and the number of interface grids in Y is 100. The number of interface grids in Z is 51 (  $4+(20-1)+(10-1)+(20-1)$  ).

The given interface model size is  $100 \times 100 \times 51$ . In fact, the largest model area we can calculate is  $(100-2 \times \text{fdx\_nghosts}) \times (100-2 \times \text{fdy\_nghosts}) \times (51-\text{fdz\_nghosts})$ . In this program, these parameters `fdx\_nghosts`, `fdy\_nghosts`, `fdz\_nghosts` are colonization 3. So the largest model area we can calculate is  $94 \times 94 \times 48$ .

(1) Example of the maximum area we can calculate without resampling:

```
number_of_total_grid_points_x = 94,
number_of_total_grid_points_y = 94,
number_of_total_grid_points_z = 48,
```

---

run\_test.sh

---

```
"layer_interp" : {
  "in_grid_layer_file" : "$EXEC_DIR/test/test_grid.gdlay",
  "refine_factor" : [ 1, 1, 1 ],
  "horizontal_start_index" : [ 3, 3 ],
  "vertical_ToFreeSurf_resample_index" : 0
}
```

---

the calculation area index range of x is 3->96,  
the calculation area index range of y is 3->96,  
the calculation area index range of z is 3->50.  
In the x direction, ghosts points is 0->2, and 97->99,  
in the y direction, ghosts points is 0->2, and 97->99,  
In the z direction, ghosts points is 0->2.

(2) Example of the partial area we can calculate without resampling:

```
number_of_total_grid_points_x = 50,
number_of_total_grid_points_y = 40,
number_of_total_grid_points_z = 25.
```

---

run\_test.sh

---

```
"layer_interp" : {
  "in_grid_layer_file" : "$EXEC_DIR/test/test_grid.gdlay",
  "refine_factor" : [ 1, 1, 1 ],
  "horizontal_start_index" : [ 20, 35 ],
  "vertical_ToFreeSurf_resample_index" : 10
}
```

---

the calculation area index range of x is 20->69,  
the calculation area index range of y is 35->74,  
the calculation area index range of z is 16->40.  
In the x direction, ghosts points is 17->19, and 70->72,  
in the y direction, ghosts points is 32->34, and 75->77,  
In the z direction, ghosts points is 13->15, and 41->42.

We provide a more complex model in the `example/prep_grid` directory.

## Chapter 2

# How to give media parameters and run wave simulations for different media types

par.json

```
"medium" : {
  "type" : "elastic_iso",
  "#code" : 1,
  "infile_layer" : "/home/usr1/testCGFD3D/prep_medium/basin.md3lay",
  "#infile_grid" : "/home/usr1/testCGFD3D/prep_medium/basin.md3grd",
  "#equivalent_medium_method" : "har",
  "#import" : "/home/usr1/testCGFD3D/output/media"
},
"is_export_media" : 1,
"media_export_dir" : "/home/usr1/testCGFD3D/output/media",

"#visco_config" : {
  "type" : "graves_Qs",
  "Qs_freq" : 1.0
}
```

- type is the flag of the solver, that is, what media do you want to simulate, and four options are supported in this code:
  - acoustic\_iso: choose this to call the solver of acoustic wave equation of isotropic media.
  - elastic\_iso: choose this to call the solver of the elastic wave equation of isotropic media.
  - elastic\_vti: choose this to call the solver of the elastic wave equation of vertical transversely isotropic media.
  - elastic\_aniso: choose this to call the solver of the elastic wave equation of anisotropic media.The media type is given in import, code, infile\_layer or infile\_grid.
- The media can be given in four ways:
  - code: Choose this flag if you want to give the media by code. You should edit **forward/md\_t.c** file and recompile the code.
  - infile\_layer: If you want to give media by the interface line, this option should be helpful. The media parameters are given by the **md3lay** file, and the file format is shown in Section 2.2
  - infile\_grid: If you want to give media by a given grid media, this option should be helpful. The media parameters are given by the **md3grd** file, the file format is shown in Section 2.1.

- `import`: If you already generated media from the above three options, and do not want to re-generate media, select this option and give the folder of the exported media.
- `equivalent_medium_method`: If you give media by `infile_layer` or `infile_grid`, equivalent medium parameterization methods can be applied. For different media, we provide different equivalent medium parameterization methods, please see Section 2.3 for detail.
- `is_export_media` should be **1** if you want to export the media, and a **`media_export_dir`** should be given.
- The `visco_config` is selected if the media is viscoelastic, the `visco_type` and `Qs` frequency are needed.

## 2.1 Format of .md3grd

If you want to get the model from the interface line, you can select the option `infile_layer`. And the media are given by a **`md3lay`** file. The file format is:

- The first line is the media type, it can be  
`one_component,`  
`acoustic_isotropic,`  
`elastic_isotropic,`  
`elastic_vti_prem,`  
`elastic_vti_thomsen,`  
`elastic_vti_cij,`  
`elastic_tti_thomsen,`  
`elastic_tti_bond,`  
`elastic_aniso_cij.`
- The second line is the number of layer (NL), if `NL > 1`, there is a designated interface, and the equivalent medium parameterization methods can be applied across the interfaces.
- The next NL lines are the number of grids in the z-direction of each layer.
- Next, the information of the interface mesh is given: `NX NY MIN_X MIN_Y SPACING_X SPACING_Y`  
`NX` and `NY` are the number of points along *x* and *y* direction;  
`MIN_X` and `MIN_Y` are the minimal *x* and *y* coordinates;  
`SPACING_X` and `SPACING_Y` are spacing between points along *x* and *y*.
- After then, the elevation, media parameters are given in every grid points. Take the `elastic_isotropic` media for example, the media parameters are read as:

```
for (nl=0; nl< NL; nl++)
  for (ip=0; ip<np[nl]; np++)
    for (iy=0; iy<NY; iy++)
      for (ix=0; ix<NX; ix++)
        fscanf(in_3grd_file, "%f %f %f %f",
              &elevation, &rho, &vp, &vs);
```

The media parameters are calculated by linear interpolation of the values at the given grid points.

For the points higher than the interface, assigned by the uppermost medium.

For different media type, different media parameters are given,

- `one_component`: var.
- `acoustic_isotropic`:  $\rho$ ,  $vp$ .
- `elastic_isotropic`:  $\rho$ ,  $vp$ ,  $vs$ .
- `elastic_vti_prem`:  $\rho$ ,  $vph$ ,  $vpv$ ,  $vsh$ ,  $vsv$ ,  $\eta$ . Please see [Dziewonski and Anderson, 1981] for detail.
- `elastic_vti_thomsen`:  $\rho$ ,  $\alpha_0$ ,  $\beta_0$ ,  $\varepsilon$ ,  $\delta$ ,  $\gamma$ . Please see [Thomsen, 1986] for detail.
- `elastic_vti_cij`:  $\rho$ ,  $c_{11}$ ,  $c_{33}$ ,  $c_{55}$ ,  $c_{66}$ ,  $c_{13}$ .
- `elastic_tti_thomsen`:  $\rho$ ,  $\alpha_0$ ,  $\beta_0$ ,  $\varepsilon$ ,  $\delta$ ,  $\gamma$ ,  $\Phi$ ,  $\theta$ . Please see [Thomsen, 1986] for detail.

- `elastic_tti_bond`:  $\rho, c_{11}, c_{33}, c_{55}, c_{66}, c_{13}, \Phi, \theta$ .  $\Phi$  is azimuth angle,  $\theta$  is the dip angle.
- `elastic_aniso_cij`:  $\rho, c_{11}, c_{12}, c_{13}, c_{14}, c_{15}, c_{16}, c_{22}, c_{23}, c_{24}, c_{25}, c_{26}, c_{33}, c_{34}, c_{35}, c_{36}, c_{44}, c_{45}, c_{46}, c_{55}, c_{56}, c_{66}$ .

if `NL > 1`, there is a designated interface; and the elevation of `ng[il]+1` needs to be the same as `ng[il]`. The equivalent medium parameterization method can be applied on the points cutting by this interface.

You can also find the example in the `example/prep_medium` directory.

## 2.2 Format of .md3lay

If you want to get the model from the interface line, you can select the option `infile_layer`. And the media are given by a `md3lay` file. The file format is:

- The first line is the media type, it can be  
`one_component,`  
`acoustic_isotropic,`  
`elastic_isotropic,`  
`elastic_vti_prem,`  
`elastic_vti_thomsen,`  
`elastic_vti_cij,`  
`elastic_tti_thomsen,`  
`elastic_tti_bond,`  
`elastic_aniso_cij.`
- The second line is the number of interface (NI).
- The third line is the information of the given interface mesh: `NX NY MIN_X MIN_Y SPACING_X SPACING_Y`  
`NX` and `NY` are the number of points along  $x$  and  $y$  direction;  
`MIN_X` and `MIN_Y` are the minimal  $x$  and  $y$  coordinates;  
`SPACING_X` and `SPACING_Y` are spacing between points along  $x$  and  $y$ .
- After then, the elevation, media parameters are given in every interface points. Take the `elastic_isotropic` media for example, the media parameters are read as:

```
for (ni=0; ni<NI; ni++)
  for (iy=0; iy<NY; iy++)
    for (ix=0; ix<NX; ix++)
      fscanf(in_3lay_file, "%f %f %f %f %f %f %f",
            &elevation,
            &rho, &rho_grad, &rho_pow,
            &vp, &vp_grad, &vp_pow,
            &vs, &vs_grad, &vs_pow);
```

`*_grad` and `*_pow` is the gradient and power of the media parameters in  $z$ -direction. The media parameters below interface are calculated by

$$var^{grid\ point} = var + (z\_interface - z^{grid\ point})^{var\_pow} * var\_grad.$$

For the points higher than the interface, assigned by the uppermost medium.

For different media type, different media parameters and the corresponding `*_grad` and `*_pow` are given, the media parameters for different media can be found in Section 2.1

You can also find examples in the `example/prep_medium` directory.



## 2.3 Note about equivalent medium parameterization methods

When there are strong interfaces in the model, using grid points to discrete models directly may cause interface error and the generation of artificial diffraction from stair step interfaces. In this code, we provide some equivalent medium parametrization methods to reduce this interface error.

If you give the model by `infile_layer` or `infile_grid`, you can apply different equivalent medium parametrization methods by give different `equivalent_medium_method`.

If the media type in the **md3grd** or **md3lay** file is `one_component`, `equivalent_medium_method` can be

- `loc`: using local value to discrete model.
- `ari`: volume integral arithmetic average.

$$\langle var \rangle = \frac{1}{\Delta V} \int_{k-1/2}^{k+1/2} \int_{j-1/2}^{j+1/2} \int_{i-1/2}^{i+1/2} var \, dx dy dz \quad (2.1)$$

- `har`: volume integral harmonic average,

$$\langle var \rangle^H = \frac{\Delta V}{\int_{k-1/2}^{k+1/2} \int_{j-1/2}^{j+1/2} \int_{i-1/2}^{i+1/2} \frac{1}{var} \, dx dy dz} \quad (2.2)$$

If the media type is `acoustic_isotropic`, `equivalent_medium_method` can be

- `loc`: using local media parameters to discrete model.
- `har`: applying harmonic average to  $\kappa$ , and applying arithmetic average to density  $\rho$ .
- `ari`: applying arithmetic average to  $\kappa$  and  $\rho$ .

If the media type is `elastic_isotropic`, `equivalent_medium_method` can be

- `loc`: using local media parameters to discrete model.
- `har`: applying harmonic average to elastic modulus, and applying arithmetic average to density  $\rho$ . Please see [Moczo et al. \[2002\]](#) and [\[Moczo et al., 2014\]](#) for detail.
- `ari`: applying arithmetic average to elastic modulus and density.

If the media type is `elastic_vti_*`, `equivalent_medium_method` can be

- `loc`: using local media parameters to discrete model.
- `har`: applying harmonic average to elastic modulus  $c_{ij}$ , and applying arithmetic average to density  $\rho$ .
- `ari`: applying arithmetic average to elastic modulus  $c_{ij}$  and density.

If the media type is `elastic_aniso_cij` or `elastic_tti_*`, `equivalent_medium_method` can be

- `loc`: using local media parameters to discrete model.
- `har`: applying harmonic average to elastic modulus  $c_{ij}$ , and applying arithmetic average to density  $\rho$ .
- `ari`: applying arithmetic average to elastic modulus  $c_{ij}$  and density.

## Chapter 3

# Source

run\_test.sh

---

```
"source_input" : {
  "in_par" : {
    "name" : "evt_by_par"
    "source" : [
      {
        "#index" : [40, 50, 30],
        "coord" : [4000, 5000, -2900],
        "wavelet_name" : "ricker",
        "ricker_center_frequency" : 0.5,
        "ricker_peak_time" : 2.0,
        "#wavelet_name" : "gaussian",
        "#gaussian_rms_width" : 2.0,
        "#gaussian_peak_time" : 0.5,
        "start_time" : 0.0,
        "end_time" : 4.0,
        "force_vector" : [1.0e13, 1.0e10, 1.0e11],
        "moment_tensor" : [1.2e13, 1.0e10, 1.0e14, 1.0e12, 1.0e11, 1.0e12]
      },
      {
        "#index" : [30, 20, 30],
        "coord" : [3000, 3000, -2900],
        "wavelet_name" : "ricker",
        "ricker_center_frequency" : 0.5,
        "ricker_peak_time" : 2.0,
        "#wavelet_name" : "gaussian",
        "#gaussian_rms_width" : 2.0,
        "#gaussian_peak_time" : 0.5,
        "start_time" : 0.0,
        "end_time" : 4.0,
        "force_vector" : [1.0e13, 1.0e10, 1.0e11],
        "moment_tensor" : [1.2e13, 1.0e10, 1.0e14, 1.0e12, 1.0e11, 1.0e12]
      }
    ]
  },
  "#in_source_file" : "$INPUTDIR/source.anasrc"    #could choose source.valsrc
},
"is_export_source" : 1,                               # if export source
"source_export_dir" : "$SOURCE_DIR",
```

---

### 3.1 JSON read source format

In the configuration of source, if `in_par` is selected, that means the source information is read by JSON. `index` is source grid index location and `coord` is source distance location, Input order is x, y, z. Select one of the two representations to determine the location of the source. `wavelet_name` include `ricker`, `ricker_deriv`, `gaussian`, `gaussian_deriv`. if `wavelet_name` choose `ricker` or `ricker_deriv`, should give `ricker_center_frequency` and `ricker_peak_time`. if `wavelet_name` choose `gaussian` or `gaussian_deriv`, should give `gaussian_rms_width` and `gaussian_peak_time`. `start_time` is source function start time, and `end_time` is source function end time. `force_vector` is single force source component, the input order is `Fx`, `Fy`, `Fz`. `moment_tentor` is double couple source component, the input order is `Mxx`, `Myy`, `Mzz`, `Myz`, `Mxz`, `Mxy`. `force_vector` and `moment_tentor` could choose together. if `in_source_file` is selected, there need to be a `.anasrc` file or `.valsrc` file, and the `.anasrc` file format is shown in Section 3.2. The `.valsrc` file format is shown in Section 3.3.

### 3.2 analytic source format

A `.anasrc` file can be given as:

---

source.anasrc

---

```

test_event_1
3 2
4.0
3000.0 2000.0 -3000.0
1.0e15 1.0e13 1.0e14
ricker
2.0 0.5
0.0
1000.0 2000.0 -3000.0
1.0e15 1.0e13 1.0e14
gaussian
2.0 0.5
1.0
1000.0 2000.0 -3000.0
1.0e15 1.0e13 1.0e14
gaussian_deriv
1.0 5.0
1.0
5000.0 2000.0 -2000.0
moment_tensor
1.0e15 1.0e13 1.0e14 1.0e15 1.0e13 1.0e14
gaussian_deriv
1.0 5.0
0.0
1500.0 2200.0 -1300.0
mechanism_angle
80.0 70.0 30.0 1.1e10 1.1 1.0e8
ricker_deriv
2.0 0.5
2.0

```

---

first line `test_event_1` is source name. second line `3 2` is force source number 3 and moment source number 2. third line `4.0` is source function time length, be equal to `end_time` subtract `start_time`. All source functions have the same time length. Next lines, the information of each source is given in order.

A force source, first line give coordinates `3000.0 2000.0 -3000.0`, input order is x, y, z. Second line give `force_vector` `1.0e15 1.0e13 1.0e14`, input order is `Fx`, `Fy`, `Fz`. Third line is wavelet name, include `ricker`, `ricker_deriv`, `gaussian`, `gaussian_deriv`. The first force source wavelet name is `ricker`. Fourth line is `ricker_center_frequency`

and ricker\_peak\_time, respectively, if wavelet name is ricker or ricker\_deriv. Fourth line is gaussian\_rms\_width and gaussian\_peak\_time, respectively, if wavelet name is gaussian or gaussian\_deriv. The first force source ricker\_center\_frequency is 2.0, ricker\_peak\_time is 0.5. The fifth line is source function start time, The first force source start time is 0.

A moment source, first line give coordinates. The first moment source coordinate is 5000.0 2000.0 -2000.0, input order is x, y, z. Second line give moment source representation type, include moment\_tensor and mechanism\_angle. The first moment source representation is moment\_tensor, so third line give monmen\_tensor component, input order is Mxx, Myy, Mzz, Myz, Mxz, Mxy. Mxx = 1.01e15, Myy = 1.0e13, Mzz = 1.0e14, Myz = 1.0e15, Mxz = 1.0e13, Mxy = 1.0e14. The fourth line is wavelet name, include ricker\_deriv, gaussian\_deriv. The first moment source wavelet name is gaussian\_deriv. The fifth line is gaussian\_rms\_width 1.0 and gaussian\_peak\_time 5.0. The sixth line is start time 0.0.

The second moment source representation type is mechanism\_angle, so next line input strike dip rake u D A. u is shear modulus, D is slip distance, unit is m, A is area. Strike = 80.0, dip = 70.0, rake = 30.0, u = 1.1e10, D = 1.1, A = 1.0e8.

### 3.3 value source format

A .valsrc file can be given as:

source.valsrc

---

```
test_event_1
1 1
0.02 20
3000.0 2000.0 -3000.0
2000.0 3000.0 -2000.0
1.0
1407.403198 14074.032227 140740.312500
5305.979492 53059.796875 530597.937500
19359.816406 193598.187500 1935981.750000
68359.867188 683598.687500 6835986.500000
233580.953125 2335809.750000 23358096.000000
772291.812500 7722918.500000 77229184.000000
2470593.750000 24705938.000000 247059376.000000
7646484.000000 76464848.000000 764648448.000000
22894096.000000 228940960.000000 2289409536.000000
66305560.000000 663055616.000000 6630556160.000000
185734304.000000 1857343104.000000 18573430784.000000
503157920.000000 5031579648.000000 50315792384.000000
1318030848.000000 13180308480.000000 131803086848.000000
3338105344.000000 33381054464.000000 333810532352.000000
8172575744.000000 81725759488.000000 817257578496.000000
19338747904.000000 193387479040.000000 1933874692096.000000
44220518400.000000 442205208576.000000 4422052085760.000000
97689878528.000000 976898818048.000000 9768988049408.000000
208449208320.000000 2084492148736.000000 20844921225216.000000
429488177152.000000 4294882099200.000000 42948817321984.000000
1.5
moment_tensor
15145.9 15145.9 15145.9 15145.9 15145.9 15145.9
28607.7 28607.7 28607.7 28607.7 28607.7 28607.7
51960.8 51960.8 51960.8 51960.8 51960.8 51960.8
90666.5 90666.5 90666.5 90666.5 90666.5 90666.5
151799.5 151799.5 151799.5 151799.5 151799.5 151799.5
243494.3 243494.3 243494.3 243494.3 243494.3 243494.3
373476.1 373476.1 373476.1 373476.1 373476.1 373476.1
546371.6 546371.6 546371.6 546371.6 546371.6 546371.6
```

```

759746.0 759746.0 759746.0 759746.0 759746.0 759746.0
999278.3 999278.3 999278.3 999278.3 999278.3 999278.3
1234194.8 1234194.8 1234194.8 1234194.8 1234194.8 1234194.8
1414798.1 1414798.1 1414798.1 1414798.1 1414798.1 1414798.1
1474371.2 1474371.2 1474371.2 1474371.2 1474371.2 1474371.2
1337506.4 1337506.4 1337506.4 1337506.4 1337506.4 1337506.4
935643.3 935643.3 935643.3 935643.3 935643.3 935643.3
228314.9 228314.9 228314.9 228314.9 228314.9 228314.9
-774273.6 -774273.6 -774273.6 -774273.6 -774273.6 -774273.6
-1994150.0 -1994150.0 -1994150.0 -1994150.0 -1994150.0 -1994150.0
-3281599.0 -3281599.0 -3281599.0 -3281599.0 -3281599.0 -3281599.0
-4430085.0 -4430085.0 -4430085.0 -4430085.0 -4430085.0 -4430085.0

```

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first line `test_event_1` is source name. second line `1 1` is force source number 1 and moment source number 1. Third line `0.02 20` is input value source time step interval 0.02 and time steps 20. Fourth line is force source coordinate `3000.0 2000.0 -3000.0`, Fifth line is moment source coordinate `2000.0 3000.0 -2000.0`. Next lines, the values of each source is given in order.

A force source, first give start time. The first force start time is `1.0`. Then give each time step force vector, each line input order is `Fx, Fy, Fz`.

A moment source, first give start time. The first moment start time is `1.5`. The second line give moment source representation type, include `moment_tensor` and `mechanism_angle`. If moment source representation is `moment_tensor`, next each line give `moment_tensor` component, input order is `Mxx, Myy, Mzz, Myz, Mxz, Mxy`. If moment source representation is `mechanism_angle`, next each line give `mechanism_angle` component, input order is `strike, dip, rake, u, D, A`.

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**Evolution of the code:**

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