# `gmd` Users Guide

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# Chapter 1

# Introduction

`gmd` is a Generalized Model Driver.

# Chapter 2

# User Input

User input is via xml control files. In general, tags use CamelCase and attributes lower case.

## 2.1 GMDSpec

`<GMDSpec>`

All input files must have as their root element `<GMDSpec>`. Recognized subelements of `<GMDSpec>` are

- `<Physics>`
- `<Permutation>`
- `<Optimization>`

Additionally, the following global elements are recognized in any element.

- `<include>`
- `<Function>`

## 2.2 Function

```
<Function id="int"
          type="str{analytic expression, piecewise linear}"
          var="str[x]" href="str" cols="list">
```

Define functions to be used elsewhere in input. `id=1` is reserved for the constant 1. function.

### 2.2.1 Examples

Analytic extression

```
<Function id="2" type="analytic expression" var="t">
  sin(t)
</Function>
```

Piecewise linear table

```
<Function id="2" type="piecewise linear">
  1 2
  2 3
  3 5
</Function>
```

Read a piecewise linear table from an external file using columns 1 and 3

```
<Function id="2" type="piecewise linear" href="./file.dat" cols="1 3"/>
```

```
% cat file.dat
# Column1 Column2 Column3
1 1 4
2 3 7
.
.
.
100 4.2 1.43
```

## 2.3 Include

```
<Include href="str"/>
```

Path to file to be included as if its contents were inplace in the input file

### 2.3.1 Example

```
<Include href="/path/to/some/file.ext"/>
```

## 2.4 Physics

```
<Physics driver="str[solid]{solid, eos}">
```

Define the physics of the simulation. Recognized subelements of `<Physics>` are

- `<Path>`
- `<Material>`
- `<Extract>`

### 2.4.1 Path

```
<Path type="str{prdef, prstate, isotherm, hugoniot}"
      format="str[default]{default, table, fcnspec}"
      nfac="int[1]" kappa="real[0]"
      tstar="real[1]" estar="real[1]" sstar="real[1]"
      amplitude="real[1]" ratfac="real[1]">
```

Define deformation paths.

#### Examples

```
<Path type="prdef" kappa="0" tstar="1" estar="-.5" amplitude="1"
      nfac="1000" ratfac="1">
  <!-- termination time, number of steps, control, Cij -->
  0 0 222222 0 0 0 0 0 0
  1 1 222222 1 0 0 0 0 0
  2 1 222222 2 0 0 0 0 0
  3 1 222222 1 0 0 0 0 0
  4 1 222222 0 0 0 0 0 0
</Path>

<Path type="prdef" nfac="100">
  0 0 444 0 0 0
  1 1 444 -7490645504 -3739707392 -3739707392
  2 1 444 -14981291008 -7479414784 -7479414784
  3 1 444 -7490645504 -3739707392 -3739707392
  4 1 444 0 0 0
</Path>
```

**Function Specification:** Function are specified as `function id[:scale]`

```
<Path type="prdef" kappa="0" tstar="1" amplitude="1" format="fcnspec"
      cfmt="222" nfac="200">
  {2 * pi} 2:1.e-1 1:0 1:0
</Path>
```

```
<Path type="prdef" format="table" cols="1 3:8" cfmt="222222" tfmt="time">
  <include href="exmpls.tbl"/>
</Path>
```

### 2.4.2 Material

```
<Material model="str">
```

Specify the material model and parameters. Material parameters are specified as individual elements.

#### Material database

```
<Matlabel db="str[MTL_PARAM_DB_FILE]">
```

Insert model parameters from a database

#### Examples

```
<Material model="elastic">
  <G>  54E+09 </G>
  <K> 124E+09 </K>
</Material>
```

```
<Material model="elastic">
  <Matlabel db="./materials.xml"> aluminum </Matlabel>
  <K> 124E+09 </K>
</Material>
```

### 2.4.3 Extract

```
<Extract format="str[ascii]{ascii, mathematic}" step="int[1]" ffmt="str[.18f]">
```

Extract variables from exodus output to different formats. Variables to be extracted are specified children of the `<Extract>` element. All components of vector and tensor variables will be extracted if only the basename is specified. Time is always extracted as the first entry of the output file.

**Examples**

Extract all components of stress and strain

```
<Extract format="ascii">
  STRESS STRAIN
</Extract>
```

Extract only the XX, YY, and ZZ components of stress

```
<Extract format="ascii">
  STRESS_XX STRESS_YY STRESS_ZZ
</Extract>
```

Extract everything

```
<Extract format="ascii">
  ALL
</Extract>
```

## 2.5 Permutation

```
<Permutation method="str[zip]{zip, combine}" seed="real[12]">
```

Permutate model input parameters to investigate sensitivities. Recognized subelements of `<Permutation>`

- `<Permutate>`

### 2.5.1 Permutate

```
<Permutate var="str"
           values="fcn{range, list, weibull, uniform, normal, percentage}"
```

Specify the paramaters to permutate. Parameters should appear elsewhere in the input file in preprocessing braces.

### 2.5.2 Example

```
<Permutation method="zip" seed="12">
  <Permutate var="K" values="weibull(125.e9, 14, 3)"/>
  <Permutate var="G" values="percentage(45.e9, 10, 3)"/>
</Permutation>
```

## 2.6 Optimization

```
<Optimization method="str[simplex]{simplex, powell, cobyla}"
              maxiter="int[25]" tolerance="real[1e-6]">
```

Optimize specified parameters against user specified objective function. Recognized subelements of `<Optimization>`

- `<Optimize>`
- `<AuxiliaryFile>`
- `<ObjectiveFunction>`

### 2.6.1 Optimize

```
<Optimize var="str" initial_value="real" bounds="list"/>
```

Specify the variable to be optimized, giving initial value and, optionally, bounds. Only the `cobyla` method accepts bounds. Variable name should occur elsewhere in the input file in braces.

### 2.6.2 AuxiliaryFile

Path to any auxiliary file needed by the optimization objective function.

```
<AuxiliaryFile href="str"/>
```

### 2.6.3 ObjectiveFunction

Path to a user defined executable script that returns the error to the optimization routine.

```
<ObjectiveFunction href="str"/>
```

An `ObjectiveFunction scriptname` is called from the command line as

```
% ./scriptname simulation_output.exo [auxiliary_file_1 [... auxiliary_file_n]]
```

### 2.6.4 Example

```
<Optimization method="simplex" maxiter="25" tolerance="1e-4" disp="0">
  <ObjectiveFunction href="opt-sig-v-time"/>
  <AuxiliaryFile href="opt-baseline.dat"/>
  <Optimize var="opt_k" initial_value="129.e9"/>
  <Optimize var="opt_g" initial_value="54.e9"/>
</Optimization>
```