MSU Internal State Variable Plasticity-Damage Model 1.0 Calibration, DMGfit Production Version Users Manual

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Summary

The Mississippi State University internal state variable (ISV) plasticity-damage model (DMG) production version 1.0 is being released along with its model calibration tool (DMGfit). The model equations are explained in Appendix A, while material model fits to sixty one (61) metals and alloys are illustrated in Appendix B. The fits were obtained by the model calibration tool DMGfit described by this document.

The DMG model is implemented as an ABAQUS user material (UMAT) subroutine for production run finite element simulations. For consistency, the same UMAT is utilized for model calibration by the DMGfit tool. The calibrated model constants can be directly merged by DMGfit into the "USER MATERIAL, CONSTANTS" section an existing ABAQUS input deck.

The DMG model UMAT and a stand-alone DMGfit tool are available, upon request, from Mark Horstemeyer <mfhorst@cavs.msstate.edu>. An online version of the DMGfit tool, along with sample experimental data, is available on the Web at http://ccg.hpc.msstate.edu/cmd.

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Quickstart Guide

On the Web, point your browser to **http://ccg.msstate.edu/cmd**, login with your credentials, select **Tools**, and then **DMGfit**. Online documentation is available by selecting **Help**. Login credentials may be requested from Tomasz Haupt <haupt@cavs.msstate.edu>. A demonstration video is available at http://ccg.hpc.msstate.edu/cmd/files/MaterialDBDemoVideo.avi.

To install DMGfit on a PC running Windows XP, request the following files from Mark Horstemeyer <mfhorst@cavs.msstate.edu>:

- 1. DMG.exe and DMG.chm the executable and user guide.
- 2. MCRinstaller.exe installer for the required runtime library.

Copy the files to, say, C:\DMG\. Run the MCRinstaller.exe, once, as a computer administrator, to install the library. Then run DMG.exe; the DMGfit GUI (see figure on the right) should appear. Select the menu item Help to display the user guide. For technical support, contact Ric Carino <rlc@cavs.msstate.edu>.

To fit the damage model to experimental data, the general strategy is to work the GUI regions in the following order: Datasets; Loading parameters; Fixed constants; Yield and yield adjustment constants; Kinematic hardening and recovery constants; Isotropic hardening and recovery constants; Torsional, Compression and Tension constants; then Damage constants. The following sequence of steps is typical:

- 1. Load all experimental datasets. For each dataset, establish the experiment settings (initial temperature, strain rate, stress units, etc), loading parameters, and fixed constants.
- 2. Start by fitting the experimental dataset with the lowest temperature and lowest rate. Temporarily exclude the rest of the datasets. If there are different tests, fit the compression datasets first, followed by tension datasets, then torsion datasets.
- 3. For the first dataset, adjust the constants as follows: yield C3; kinematic hardening C9 and recovery C7; isotropic hardening C15 and recovery C13.
- 4. Restore second dataset. If it has a different temperature than the first, adjust the constants as follows: {C3, C4} if yield is temperature dependent; then {C10, C8, C16, C14}. If the dataset has a different strain rate, adjust C1 and C5 if yield is strain rate dependent, then {C9, C7, C11} and {C15, C13, C17}.
- 5. Repeat step 4 for the rest of the datasets. If adjusting the temperature dependence constants (even Cs) does not produce good models for high temperature data, adjust C19 and C20. Adjust torsion, compression and tension differentiation constants, if adding stress state dependent experimental data.
- 6. Adjust damage constants. To track the evolution of the damage state variable, select *Display | Plot a state variable | Damage*. Readjust constants in other boxes as necessary.
- 7. Record your results. Select *File | Write | A binary restart file* to create a record of your session, to resume it later. Select *File | Write | A "USER MATERIAL, CONSTANTS" for ABAQUS* to create a text file containing the properties, suitable for inclusion in an ABAQUS input deck. Select *File | Write | A "fitted constants+constrains" file* to create a text file record of the fitted constants, which may be used as starting values in future fitting sessions.

Introduction

The Mississippi State University internal state variable (ISV) plasticity-damage model (DMG) production version 1.0 is based from the ISV plasticity formulation of Bammann [1990] with the addition of porosity [Bammann et al., 1993] and the void nucleation, growth, and coalescence rate equations that admit heterogeneous microstructures [Horstemeyer et al, 2000]. The model is implemented as an ABAQUS user material subroutine (UMAT). The model calibration routine DMGfit was developed by Cariño [2007], which updated the original BFIT routine by Lathrop [1996]. The calibrated model constants can be directly merged into the "USER MATERIAL, CONSTANTS" section an existing ABAQUS input deck.

The pre-production release of DMGfit is based on a C translation of an earlier version of the DMG code (Research version 0.9 UMAT in Figure 1 below), instead of the state-of-the-art version used in production-run finite element simulations. Synchronizing DMGfit with the latest version of the ABAQUS UMAT implementation of the damage model by revising the C translation is a tedious process, a duplication of efforts and an error-prone procedure. Further, this revision has to be performed for future extensions to the damage model since these extensions are tested by MSU researchers on the UMAT implementation. In order to avoid these inefficiencies in the software maintenance process and to keep DMGfit always synchronized with the current version of the damage model, DMGfit has been modified to invoke the damage model UMAT.



Figure 1. Development roadmap for DMG and DMGfit. The earlier DMGfit uses an outdated C-translation of DMG UMAT. DMGfit was upgraded to link with the production DMG UMAT.

Input-Process-Output

Production version 1.0 of the DMG model is specified by 54 constants or material properties. Some of these constants, like bulk modulus, shear modulus, and melting temperature, are fixed for a given material and are obtained from the literature. Microstructure constants, such as average radius of voids, average size of particles, and average grain size, are calculated from microscope images of samples of the material. The remaining constants are fitted, using DMGfit, from stress-strain experiments on the samples.

DMGfit expects the measurements associated with a sample of the material to be recorded in a "DMGfit data set". This data set contains experimental stress-strain data, fixed constants for the material, fixed constants for the specific sample, and the loading parameters to simulate the experiment. DMGfit solves an optimization problem to calculate the model constants that are not included in a data set.

Figure 2 is a high-level description of the process to calibrate the damage model using DMGfit. Three codes are involved: an optimization routine (in MATLAB), a material point simulator (in Fortran), and the damage model UMAT (in Fortran). Collaborators may request these codes from Mark Horstemeyer <mfhorst@cavs.msstate.edu>. The inputs to the process are estimated initial values for the constants and experimental stress-strain data. Candidate model constants are generated by the optimization routine. These constants are passed as material properties to the Material Point Simulator. The simulator invokes the damage model UMAT with the properties to produce a model stress-strain curve. The process successfully terminates when the model curve gives a good correlation to the experimental data.



Figure 2. Overview of model calibration process using DMGfit.

Graphical User Interface

The remainder of this report describes the user interface of the stand-alone version of DMGfit. The documentation for the Web version of DMGfit is online at http://ccg.hpc.msstate.edu/ccgportlets/apps/cmd/html/help/Help.htm.

A snapshot of the DMGfit GUI in operation, annotated to highlight the logical groupings of the controls, is shown by Figure 3.



Figure 3. The logical groupings of controls in the DMGfit GUI.

Brief descriptions of the controls are:

- **Menu items** file operations, optimization options, plot appearance settings, and help links. **Data set controls** - describes the experimental stress-strain data measured from a sample of the material under consideration.
- **Loading parameters** parameters for the material point simulator to model the experiment. **Fixed constants** - damage model constants that are fixed, from the literature, or measured

from microscope images of the sample represented by the data set.

Single parameter study controls - activated by a right-click on a fitted constant, to investigate the *solitary* effect of the constant on the model.

Parameter set controls - operations that apply to the current *set* of fitted constants. **Fitted constants** - damage model constants to be calibrated.

The controls are further described in the sections that follow.

Menu Items

DMGfit for Damage Model v.1.0 File Optimization Display Help

File

Load

- A data file load a DMGfit data set file (see Data set controls).
- A ''fixed constants'' file load a text file containing lines with the format "name=value", where "name" is a fixed constant in the damage model
- Stress-strain from Excel file load stress-strain data from an Excel file (see Stressstrain data).
- A "fitted constants+constraints" file load a text file containing lines with the format "name=value fix min max", where "name" is a fitted constant in the damage model, fix is 0 or 1 to indicate that the value is to be fixed during optimization, and [min, max] is the range of permissible values for the constant if fix=0.
- A "fitted constants only" file load a text file containing lines with the format "name=value", where "name" is a fitted constant in the damage model
- A binary restart file a record of a previous fitting session.

Write

- A post-processing info text file create a text file containing the model constants, data points and model points for post-processing, most likely to produce publication-quality plots. This file can be imported as comma-separated-values (CSV) into a spreadsheet program like MS Excel.
- A Gnuplot script for the plots create a Gnuplot script for the plots and the parameter set on focus.
- A binary restart file create a record of the state of a fitting session. Loading the file with Load | Binary restart file continues the session from the point the file was saved.
- A data file create a template for a data file from a loaded data set.
- A "fitted constants+constraints" file create a file for the fitted constants and their constraints, from the parameter set on focus.
- A "USER MATERIAL, CONSTANTS" for ABAQUS write the fixed constants and fitted constants into an existing ABAQUS input deck, or create such a section if the deck does not exist yet.

Utilities

- Extract constants from ABAQUS input extract the "USER MATERIAL, CONSTANTS=54" section of an ABAQUS input deck, then write the constants into a text file that can be retrieved using Load | A "fitted constants only" file or Load | A "fixed constants" file
- **Extract constants from legacy param file** extract the constants from a BFIT param file, then write the constants into a text file that can be retrieved using **Load** | **A "fitted constants only" file**

Extract constants from 1-column props file - read a text file with a single column of 54 numbers, then write the numbers as damage model constants into a text file that can be retrieved using Load | A ''fitted constants only'' file or Load | A ''fixed constants'' file

Print - open a print preview of the DMGfit window.

Reset - re-initialize DMGfit.

Replay - execute a replay file, which is a text record of the actions in a previous DMGfit session.

Quit - write a replay file for the session if wanted; then quit DMGfit.

Optimization

Use fmincon - use MATLAB's fmincon(), after setting some options.
Use fminsearch - use MATLAB's fminsearch(), after setting some options (the default).
Use lsqnonlin - use MATLAB's lsqnonlin(), after setting some options.
Use patternsearch - use MATLAB's patternsearch(), after setting some options.
Use genetic algorithm - use MATLAB's ga(), after setting some options.
Estimate C3 and C4 - calculate C3 and C4 from two datasets with different temperatures.

Display

Plot settings - open a dialog to modify the appearance of the plot area.
Plot a state variable - open a dialog to select the state variable to plot.
Toggle Hide|Show model plots - enable or disable plotting of the model curves.
Font size - set the font size; fonts in the plot area will be changed when the plot is updated.
Font name - set the font style; fonts in the plot area will be changed when the plot is updated.

Help

Help topics - display the help file.

About - display the current version of DMGfit and the contact information of the developer/maintainer.

Data Set Controls

A DMGfit data set is a file that contains experimental stress-strain data, descriptive elements of the experimental sample, the loading parameters to model the stress-strain data, and the fixed constants associated with the sample. A data set file is retrieved from disk using **File | Load | A data file**. The information contained in the file are automatically displayed when the data set is selected using the **DATA SET** popup.

DATA SET	F AE42-150	c	•			
Average	Reduce	Remove	Exclude			
Material	AE42					
Source	David Oglesby					
Comment	Strain rate n	ot given				
Tag	AE42-150c					
Test type	1	Units	5			
Tinit	423.15	Rate	0.001			

Hovering the mouse pointer over a control causes a pop-up to appear. For examples, Average experiments on Average; and Stress units: 1=psi 2=Ksi 3=Pa 4=KPa 5=MPa 6=GPa on Units 5

The stress-strain data is described by the following elements:

Material - a descriptive name for the material.
Source - who provided the experimental data.
Comment - any remarks about the data or the experiment.
Tag - a tag to differentiate data sets.
Test type - 1=Tension, 2=Compression, 3=Torsion, 4=Tension+Compression, 5=Compression+Tension
Units - stress units: 1=psi 2=Ksi 3=Pa 4=KPa 5=MPa 6=GPa
Tinit - initial temperature.
Rate - strain rate.

Except for the stress-strain data which are plotted, the above elements may be edited directly. Changes will take effect after clicking Apply changes (see **Parameter set controls**).

Operations on a data set are:

Average starts a dialog to average two or more experiments with the same Test type, Tinit and Rate. A dialog ensues to select the data sets to be averaged, the number of strain values (say, *npoints*) to be generated, and a file name for the differences between the average stress and experimental stresses. A "new" data set will be created from the selected experiments and given an "Averaged" Tag. The stress values will be the average of the interpolated stress at *npoints* equally-spaced values between the minimum and maximum experimental strain values. The descriptive elements, loading parameters and fixed constants are derived from the first data set that was selected. If a file name was specified, the file will be created with the following columns: the strain values, the averaged interpolated stress, the differences between the average stress and experimental stresses, and the minimum and maximum differences

Reduce disables all controls and makes visible controls for reducing the number of stressstrain pairs in the data set. Thousands of data points translate into long running times for the fitting procedure since an estimate of the error at each point has to be calculated.



The **left slider** moves the left endpoint, while the **right slider** moves the right endpoint. The **Fitter** button invokes a filtering routine to find points such that a line drawn between the points approximates the data between those points to a relative error not exceeding the filter tolerance. For comparison, the points found are superimposed on the original dataset. <u>Cancel</u> discards changes to the data set, deactivates the reduction controls, and enables the rest of the controls. <u>Accept</u> creates a new data set from the filtered/reduced stress-strain data points of the data set on focus, prefixing "reduced-" to the tag. The data set on focus is automatically excluded, and the reduced data set becomes the focus.

Remove deletes the data set on focus from memory.

Exclude inhibits the data set from being plotted and fitted. The data set Tag is annotated with "(excl)". If an excluded data set is selected for focus, this button changes to Restore, for the opposite effect.

Tips:

 If the data set on focus has Test type=2 (compression) and the stress-strain data are negative, it is necessary to "flip" the data so that the strain and stress values become positive. A rightclick on Average exposes the context menu

AVE	Pra	nel	Reduce	Remove
Mε		Flip	compressio	on data set
c	. 1	•		

for this purpose.

2. A right-click on Remove exposes the context menu

Remove Exclude Remove some data sets

to remove two or more data sets. A dialog ensues, to select the data sets to be removed.

3. A right-click on **Exclude** exposes the context menu

Exclude Exclude some data sets

to exclude two or more data sets. A dialog ensues, to select the data sets to be excluded or restored.

4. A right-click on the value of Tinit exposes the context menu



to enter the initial temperature in Celsius or Fahrenheit.

Loading Parameters

A set of loading parameters is associated with each experimental data. These parameters are used by the material point simulator to model the data. The values of these parameters are automatically displayed upon

Loading parameters							
Period	234.272	e-load	0				
Incr	100	Angle	0.51				

Descriptions of the loading parameters:

Period - simulation time period; the final simulated strain value will be (Period*Rate)
e-load - for Test type=4 or 5; the strain value where the load direction changes; otherwise 0.
Incr - number of strain increments (up to 1000) for the simulation; the execution time of the simulation is directly proportional to this value.

Angle - the bi-axial loading ratio.

Tips:

- 1. If Period is 0, a model will not be generated for the data set.
- 2. If a model curve is a "zigzag" or is "wavy", increase the corresponding Incr parameter.
- 3. A right-click on Period exposes the context menu



to automatically calculate all periods so that all data models will end at a common strain value, or at the respective final strain values. For example, in the figure below, the Periods were automatically computed to model up to strain=0.02.



Fixed Constants

Some constants in the damage model are fixed for the material. These are obtained from the literature or measured from microscope images of the sample used in the experiment. The **Fixed constants** associated with a sample are automatically displayed upon selection of the data set. These values may be edited directly. Changes will take effect after clicking Apply changes (see **Parameter set controls**). Hovering the mouse pointer

over a value causes a pop-up describing the constant to appear.

Descriptions of the fixed constants:

G - Shear modulus (default: 0)

Bulk - Bulk modulus (default: 0)

a - Constant to adjust shear modulus due to variation of temperature from initial (default: 0)

b - Constant to adjust bulk modulus due to variation of temperature from initial (default: 0)

Tmelt - Melt temperature (K) (default: 3000)

heat - Heat generation term (default: 0.34)

Kic - Particle fracture toughness (default: 1000)

- nv Particle McClintock damage constant (default: 0.3)
- dn Particle average size (default: 0)
- fn Particles volume fraction, must be greater than 0 (default: 1e-3)
- dcs0 Reference grain size or dendrite cell size, must be greater than 0 (default: 30)

dcs - Grain size or dendrite cell size of experiment sample (default: 30)

r0 - Initial radius of a spherical void (default: 0)

volF - Initial void volume fraction (default: 0)

Tips:

- 1. The constants fn and dcs0 should not be zero; these are divisors inside the UMAT.
- 2. If r0 and volF are zero, the damage constants will have no effect on the model.
- 3. If G or Bulk is 0, a model will not be generated for the data set.
- 4. A right-click on G or Bulk exposes a context menu

G 16667 Bulk 50000 %

a 🔽 💿 (G, Bulk) from (Youngs, Poisson)

to calculate the bulk and shear moduli from the Young modulus and Poisson ratio if these are the known quantities.

Fixed constants

G	16667	Bulk	50000
а	0	b	0
Tmelt	950	heat	0.32
nv	0.3	Kic	1000
dn	0	fn	0.001
dcs	30	dcs0	30
rO	0	volF	0

Single Parameter Study Controls

Parameter study on C3					
0	145	1e+006			
Min		▶ Max			
Cancel	Scale=10^ 0	Acce	pt		

A right-click on the current value of a fitted constant activates the parameter study controls and disables the rest of the DMGfit controls. In the above example, the right-click was on C3 145

The range constraint for C3 is transferred to the **Min** and **Max** edit fields as well as to the slider limits. The current value is transferred to the **Current** edit field and is used to initially position the slider.

The **Min**, **Current** and **Max** values may be edited directly. Changes will take effect after clicking anywhere in the window outside of the edit boxes.

Operating the slider decreases or increases the **Current** value, depending on where the slider bar is positioned. The model plots are updated automatically with each change of the current value.

If the **Min** value is close to the **Max** value, the slider may not work properly due to the resulting low resolution of the slider steps. In this case, setting the scale (Scale=10[^] edit box) to a negative integer may restore the expected functionality. Similarly, for a large range of values, setting the scale to a positive integer may prevent huge jumps in **Current** when the slider is operated.

Accept creates a new parameter set from the set currently on focus, with the current value and constraint for the selected constant coming from the parameter study controls. This also exits the parameter study. **Cancel** exits the parameter study without any changes to the constant.

Parameter Set Controls

Optimize parameters	Apply changes	
PARAMETER SET 44 of 44		Remove

DMGfit keeps track of the sets of fitted constants (i.e., parameter sets) attempted during a session. The **PARAMETER SET slider** selects the set which will be used to generate the model plots.

Remove discards the current parameter set. The next available parameter set becomes the focus.

Optimize parameters invokes the optimization routine to automatically update the values in the parameter set.

Apply changes commits changes to any displayed value (experiment descriptions, loading parameters, fixed constants, fitted constants) or check box in the GUI. If fitted constants are changed, a new parameter set is created from the displayed values. The new set automatically becomes the parameter set on focus.

Fitted Constants

Yield a	Yield and yield adjustment constants Kinematic hardening and recovery constants							Isotrop	ic hardening	g and r	ecovery	/ constants					
C3	145		C4	0	-	C9	4874.6		C10	3.71		C15	1079		C16	0.021	
C1	0	•	C2	0		C7	3.53		C8	103.067		C13	2.05284		C14	1000	
C5	1e-005	•	C6	0		C11	0		C12	0	•	C17	0	•	C18	0	<u>-</u>
C19	0.0065		C20	475								Dama	ge constant	s			
Long ra	ange hardeni	ing an	d recov	ery constan	ıts	Torsio	nal, Compre	ssion	an	0		Ccoef	0	V	NTD	0	•
C23	2.677		C24	8.48		äle	nsion const	ants	bn	0		cd1	0	v	CTD	0	v
C21	0.668		C22	591.5		Са	0	•	cn	0		cd2	0	•	CAcon	0.5	•
C25	0	•	C26	0		Cb	0		Z	0							

For each fitted constant, there are three controls: the **label** static text, the current **value** edit box, and the **fix** check box.

For example, consider $\boxed{^{C3}$ $\boxed{^{145}}$ $\boxed{\mathbf{v}}$. Hovering the mouse pointer on the edit box will activate a popup that displays a description and current range constraint for the constant.

Constant term in Y(T) Arrhenius-type equation which is the rate independent yield stre	ss
Min= 0	
Max= 1e+006	
Fix=1	

The current value can be changed directly. The change will take effect when Apply changes is clicked.

A right-click on the current value activates the **Single parameter study controls** and disables the rest of the DMGfit controls.

If the check box is unchecked, the constant may be updated when Optimize parameters is clicked. If checked, the constant will be fixed (not changed) by the optimization routine.

Tips:

- 1. Leave C5 to the default nonzero value if yield is NOT strain rate dependent. C5=0 does not produce a good model.
- 2. Several values can be modified and/or fixed before Apply changes is clicked. A new parameter set is created from the set on focus, using the updated values. The new set automatically becomes the focus.
- 3. Depending on the optimization routine, a constant set to zero may be respected, and may remain zero *even if it is not fixed*. For the constant to participate in fitting, set it initially to a "small" non-zero value.
- 4. The group ordering of the fitted constants in the GUI suggests the fitting strategy. After establishing the fixed constants, the suggested order for the constants to be fitted is:
 - 1. Yield and yield adjustment constants,
 - 2. Kinematic hardening and recovery constants,
 - 3. Isotropic hardening and recovery constants,
 - 4. Long range hardening and recovery constants,
 - 5. Torsional, Compression & Tension differentiation constants, and
 - 6. Damage constants.

Descriptions of the constants:

- C1 Constant term in V(T) Arrhenius-type equation which determines the magnitude of rate dependence on yielding (units=MPa)
- C2 Temperature dependent activation term in V(T) Arrhenius-type equation (units=Kelvin)
- C3 Constant term in Y(T) Arrhenius-type equation which is the rate independent yield stress (units=MPa)
- C4 Temperature dependent activation term in Y(T) Arrhenius-type equation (units=Kelvin)
- C5 Constant term in f(T) Arrhenius-type equation which determines the transition strain rate from rate independent to dependent yield (units=1/sec)
- C6 Temperature dependent activation term in f(T) Arrhenius-type equation (units=Kelvin)
- C7 Constant term in rd1 equation which describes the kinematic dynamic recovery (units=1/MPa)
- C8 Temperature dependent activation term in rd1 equation (units=Kelvin)
- C9 Constant term in h1 equation which describes the kinematic anisotropic hardening modulus (units=MPa)
- C10 v term in h1 equation (units=MPa/Kelvin)
- C11 Constant term in rs1 equation which describes the kinematic static recovery (units=1/(MPa sec))
- C12 Temperature dependent activation term in rs1 equation (units=Kelvin)
- C13 Constant term in rd2 equation which describes the isotropic dynamic recovery (units=1/MPa)
- C14 Temperature dependent activation term in rd2 equation (units=Kelvin)
- C15 Constant term in h2 equation which describes the isotropic hardening modulus (units=MPa)
- C16 Temperature dependent activation term in h2 equation (units=MPa/Kelvin)
- C17 Constant term in rs2 equation which describes the isotropic static recovery (units=MPa/sec)
- C18 Temperature dependent activation term in rs2 equation (units=Kelvin)
- C19 Multiplication term in equation 1 + tanh(c19(c20-T)), which is an adjustment to the yield strength over a large temperature range (units=1/Kelvin)
- C20 Used in yield strength adjustment equation (units=Kelvin)
- C21 Long range kinematic dynamic recovery rd3 (units=1/MPa)
- C22 Temperature dependent activation in rd3 dynamic recovery equation (units=Kelvin)
- C23 Long range kinematic hardening h3 (units=MPa)
- C24 Temperature dependent activation in h3 equation (units=MPa/Kelvin)
- C25 Long range kinematic static recovery rs3 (units=1/(MPa sec))
- C26 Temperature dependent activation in rs3 static recovery equation (units=Kelvin)
- Ca Torsional softening constant in kinematic and isotropic hardening and dynamic recovery equations
- Cb Tension/compression asymmetry constant in kinematic and isotropic hardening and dynamic recovery equations
- an Torsional constant a in void-crack nucleation model
- bn Tension/compression constant b in void-crack nucleation model
- cn Triaxiality constant c in void-crack nucleation model
- Z Grain size or dendrite cell size exponent
- Ccoef Coefficient constant in void-crack nucleation model

cd1 – Void impingement coalescence factor, D=nucleation*void volume*coal
cd2 – Void sheet coalescence factor, D=nucleation*void volume*coal
NTD – Void-crack nucleation Temperature Dependence
CTD – Void coalescence Temperature Dependence
CAcon - Cocks Ashby void growth constant

Stress-strain Data

For compatibility with the original BFIT, text files with the following formats for stress-strain data are recognized:

The other elements of the data set (Tag, Source, Comment, Initial temperature, Strain rate, etc) may be entered using the **Data set controls**.

DMGfit is able to load stress-strain data from Excel files. Use the menu item **File** | **Load data from Excel file** to open an MS Excel file and read a rectangular array of numbers. Column indices for the strain and stress values and other data set elements are specified in the subsequent dialog.

Use the menu item **File** | **Write a data file** to create a template for data files from a loaded data set. Data files so created may later be loaded using **File** | **Load a data file**.

References

- Bammann, D. J., "Modeling Temperature and Strain Rate Dependent Large of Metals," *Applied Mechanics Reviews*, Vol. 43, No. 5, Part 2, May, 1990.
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- Lathrop, J.F., (Dec 1996). BFIT—A Program to Analyze and Fit the BCJ Model Parameters to Experimental Data Tutorial and User's Guide. SANDIA REPORT SAND97-8218. UC-405, Unlimited Release, Printed December 1996.

Appendix A. MSU ISV DMG 1.0 Production Model Equations

The MSU ISV DMG 1.0 production material model is given by the following equations. The pertinent equations in this model are denoted by the rate of change of the observable and internal state variables. The equations used within the context of the finite element method are given by,

$$\overset{\circ}{\underline{\sigma}} = \underline{\dot{\sigma}} - \underline{W}^{e} \underline{\sigma} - \underline{\sigma} \underline{W}^{e} = \lambda (1 - D) tr(\underline{D}^{e}) \underline{I} + 2\mu (1 - D) \underline{D}^{e} - \frac{\dot{D}}{1 - D} \underline{\sigma}$$
Equation A.1
$$D^{e} = D - D^{in}$$
Equation A.2

$$\underline{D}^{in} = f(T) \sinh \left[\frac{\|\underline{\sigma}' - \underline{\alpha}\| - \{R + Y(T)\}\{1 - D\}}{V(T)\{1 - D\}} \right] \frac{\underline{\sigma}' - \underline{\alpha}}{\|\underline{\sigma}' - \underline{\alpha}\|}$$
Equation A.3

$$\frac{\overset{o}{\alpha}}{\underline{\alpha}} = \underline{\dot{\alpha}} - \underline{W}^{e} \underline{\alpha} + \underline{\alpha} \underline{W}^{e} = \left\{ h(T) \underline{D}^{in} - \left[\sqrt{\frac{2}{3}} r_{d}(T) \| \underline{D}^{in} \| + r_{s}(T) \right] \| \underline{\alpha} \| \underline{\alpha} \right\} \left[\frac{DCS_{0}}{DCS} \right]^{z}$$
Equation A.4

$$\dot{R} = \left\{ H(T)\underline{D}^{in} - \left\lfloor \sqrt{\frac{2}{3}}R_d(T) \right\| \underline{D}^{in} \right\| + R_s(T) \left\| R^2 \right\} \left[\frac{DCS_0}{DCS} \right]^z$$
Equation A.5
$$\dot{D} = \left[\dot{\phi}_{particles} + \dot{\phi}_{pores} \right] c + \left[\phi_{particles} + \phi_{pores} \right] \dot{c} ,$$
Equation A.6

$$\dot{\phi}_{\text{particles}} = \dot{\eta}v + \eta\dot{v}$$

$$\dot{\eta} = \left\| \underline{D}^{in} \right\| \frac{d^{\frac{1}{2}}}{K_{IC} f^{\frac{1}{3}}} \eta \left\{ a \left[\frac{4}{27} - \frac{J_3^2}{J_2^3} \right] + b \frac{J_3}{J_2^{\frac{3}{2}}} + c \left\| \frac{I_1}{\sqrt{J_2}} \right\| \right\} \exp \left(-\frac{C_{\eta T}}{T} \right)$$
Equation A.8
3 $\left[3 V(T) \sigma_{T} - \left(-\frac{V(T)}{T} \right) \right] = c \left[\frac{1}{\sqrt{J_2}} + \frac{1}{\sqrt{J_2}} \right]^{\frac{V(T)}{V(T)}} = c$

Equation A.7

$$\dot{v} = \frac{3}{2} v \left[\frac{3}{2} \frac{V(T)}{Y(T)} \frac{\sigma_H}{\sigma_{vm}} + \left(1 - \frac{V(T)}{Y(T)} \right) (1 + 0.4319) \right]^{-\gamma_V(T)} \underline{D}^{in}$$
Equation A.9

$$\dot{c} = C_{coal} \left[\eta \dot{v} + \dot{\eta} v \right] \exp \left(C_{CT} T \right) \left(\frac{DCS_0}{DCS} \right)^z$$
Equation A.10

$$\dot{\phi}_{pores} = \left[\frac{1}{\left(1 - \phi_{pores}\right)^{m}} - \left(1 - \phi_{pores}\right)\right] \sinh \left\{\frac{2\left(2^{V(T)}/Y(T)^{-1}\right)}{\left(2^{V(T)}/Y(T)^{+1}\right)} \frac{\sigma_{H}}{\sigma_{vm}}\right\} \left\|\underline{D}^{in}\right\|$$
Equation A.11

The rate equations are generally written as objective rates $(\overset{o}{\sigma}, \overset{o}{\alpha})$ with indifference to the continuum frame of reference assuming a Jaumann rate in which the continuum spin equals the elastic spin ($\underline{W} = \underline{W}^e$). The internal state variable (ISV) equations are functions of the observable variables (temperature, stress state, and rate of deformation). In general, the rate equations of generalized displacements, or thermodynamics fluxes, describing the rate of change may be written as independent equations for each ISV or as derivatives of a suitably chosen potential function arising from the hypothesis of generalized normality. An advantage of assuming generalized normality, although somewhat restrictive, is unconditional satisfaction of the Kelvin inequality of the second law of thermodynamics (nonnegative intrinsic dissipation), i.e.

$$\underline{\sigma}: \underline{D}^{in} - \underline{b}: \underline{\alpha} - \kappa \bullet \dot{R} - \phi \bullet \dot{D} \ge 0.$$
 Equation A.12

The selection of the ISVs may, in principle, be somewhat arbitrary, but the kinematic hardening, isotropic hardening, and damage rate equations are physically motivated and strongly influence the history of the material. The ISV model accounts for deviatoric inelastic deformation resulting from the presence of dislocations in crystallographic material, dilatational deformation, and ensuing failure from damage progression. Damage will reduce the material strength, enhance the inelastic flow, and soften the elastic moduli.

In Equation A.1, the elastic Lame constants are denoted by λ and μ . The elastic rate of deformation (\underline{D}^e) results when the flow rule as shown in Equation A.2 is subtracted from the total deformation (\underline{D}), which is defined by the boundary conditions.

The independent variables for the inelastic rate of deformation are given in Equation A.3 as the stress, temperature, and internal state variables. This is similar to power law and Garofalo equations for creep except that the ISVs are now included. The deviatoric inelastic flow rule, \underline{D}^{in} , encompasses the regimes of creep and plasticity and is a function of the temperature, the kinematic hardening internal state variable ($\underline{\alpha}$), the isotropic hardening internal state variable (R), the volume fraction of damaged material (D), and the functions f(T), V(T), and Y(T), which are related to yielding with Arrhenius-type temperature dependence. The function Y(T) is the rate-independent yield stress. The function f(T) determines when the rate-dependence affects initial yielding. The function V(T) determines the magnitude of rate-dependence on yielding. These functions are determined from simple isothermal compression tests with different strain rates and temperatures,

$$V(T) = C_1 \exp\left(\frac{-C_2}{T}\right), \ Y(T) = C_3 \exp\left(\frac{C_4}{T}\right), \ f(T) = C_5 \exp\left(\frac{-C_6}{T}\right).$$
 Equation A.13

The kinematic hardening internal state variable, $\underline{\alpha}$, reflects the effect of anisotropic dislocation density, and the isotropic hardening internal state variable R, reflects the effect of the global dislocation density. As such, the hardenings are cast in a hardening-recovery format that includes dynamic and static recovery. The functions $r_s(T)$ and $R_s(T)$ are scalar in nature and describe the diffusion-controlled static or thermal recovery, while $r_d(T)$ and $R_d(T)$ are scalar functions describing dynamic recovery. Hence, the two main types of recovery that are exhibited by populations of dislocations within crystallographic materials are captured in the ISVs. The anisotropic hardening modulus is h(T), and the isotropic hardening modulus is H(T).

The hardening moduli and dynamic recovery functions account for deformation-induced anisotropy arising from texture and dislocation substructures by means of stress-dependent variables. Miller *et al.* (1995) showed that by using J_3 in the hardening equations the different hardening rates between axisymmetric compression and torsion (torsional softening) were accurately captured. Miller *et al.* (1995) and Horstemeyer *et al.* (1995) included this feature in

the Bammann ISV model as

$$r_{d}(T) = C_{7} \left(1 + C_{19} \left[\frac{4}{27} - \frac{J_{3}^{2}}{J_{2}^{3}} \right] \right) \exp\left(-C_{8} / T \right),$$
 Equation A.14
$$h(T) = \left\{ C_{9} \left(1 + C_{99} \left[\frac{4}{27} - \frac{J_{3}^{2}}{J_{2}^{3}} \right] \right) \right\} - C_{9} T,$$
 Equation A.15

$$r_{s}(T) = C_{11} \exp\left(\frac{-C_{12}}{T}\right),$$
 Equation A.16

$$R_{d}(T) = C_{13}\left(1 + C_{21}\left[\frac{4}{27} - \frac{J_{3}^{2}}{J_{2}^{3}}\right]\right) \exp\left(-C_{14}/T\right),$$
 Equation A.17

$$H = \left\{ C_{15} \left[1 + C_{22} \left[\frac{4}{27} - \frac{J_3^{'2}}{J_2^{'3}} \right] \right] \right\} - C_{16}T,$$
 Equation A.18

$$R_{s}(T) = C_{17} \exp\left(\frac{-C_{18}}{T}\right),$$
 Equation A.19

where $J_2 = \frac{1}{2} (\underline{\sigma} - \underline{\alpha})^2$ and $J_3 = \frac{1}{3} (\underline{\sigma} - \underline{\alpha})^3$. The deviatoric stress $\underline{\sigma}'$ is expressed in indicial notation as

$$\sigma_{ij} = \sigma_{ij} - \frac{1}{3}\sigma_{ii}$$
. Equation A.20

The damage variable D represents the damage fraction of material within a continuum element. The mechanical properties of a material depend upon the amount and type of microdefects within its structure. Deformation changes these microdefects, and when the number of microdefects accumulates, damage is said to have grown. The notion of a damaged state in continuum field theory emerged when Kachanov (1958) introduced a damage variable to describe the microdefect density locally in a creeping material. The idea was that damage could be measured by the volume fraction of voids under creep conditions. Rabotnov (1969) furthered this notion with a rate equation of void density.

Equation A.6 introduces the void volume fraction (porosity) as damage. By including damage, D, as an ISV, different forms of damage rules can easily be incorporated into the constitutive framework. Bammann *et al.* (1993; 1996) and Horstemeyer (1992, 1993) have demonstrated the applicability of the Cocks and Ashby (1980) void growth rule used as the damage rate equation in the ISV model.

The generalized thermodynamic force conjugate, ϕ , is often referred to as the energy release rate for elastic brittle materials and the J-integral for inelasticity. In essence, an increment of damage will have associated energy released per unit damage extension as new damaged area (or volume) is developed.

The damage progression is divided into void nucleation and growth from silicon particles and

from pores. Coalescence is introduced to reflect pore-pore interactions and silicon-pore interactions as expressed in Equation A.10. The void nucleation evolution described by Equation A.8 is discussed in length by Horstemeyer and Gokhale (1999). The void growth related to silicon particles, Equation A.9, is that from Budiansky *et al.* (1982). Other forms can be used and evaluated (cf., Horstemeyer *et al.*, 1999), but this equation allows for a strain rate sensitivity in relation the plasticity model (m=V(T)/Y(T)). For the porosity evolution, the Cocks and Ashby (1981) void growth rule is used as shown in Equation A.11.

Appendix B. MSU ISV DMG 1.0 Production Version: Material Model Fits

B1 A1 A356 T6 allow	B2 A1 90% pure allow
$\begin{array}{c} \text{D1. Al } \text{A550-10 alloy} \\ \text{D2. Al } 1000 \text{ T0 alloy} \\ \end{array}$	D2. Al 99% pute anoy
B5. Al 1000-10 alloy	D4. AI 1100 alloy
B5. AI 2024-10 alloy	B6. AI 2024 1351 alloy
B7. Al 2024-14 alloy	B8. AI 5083 alloy
B9. Al 5086-T6 alloy	B10. Al 6022 alloy
B11. Al 6061-T0 alloy	B12. Al 6061-T6 alloy
B13. Al 6061-T651 alloy	B14. Al 7039 alloy
B15. Al 7075-T0 alloy	B16. Al 7075-T6 alloy
B17. Al 7075-T65l alloy	B18. 300mmarsteel alloy
B19. 1006 steel alloy	B20. 1020 steel alloy
B21. 10b22 steel alloy	B22. 304L SS alloy
B23. 321 SS alloy	B24. 410 SS alloy
B25. 4140 steel alloy	B26. 4340 steel alloy
B27. A286 steel alloy	B28. AF steel alloy
B29. Armco Iron	B30. Armor Steel
B31. C1008 Steel	B32. Capeleciron
B33. FC0205 steel alloy	B34. HY80 steel alloy
B35. HY100 steel alloy	B36. HY130 steel alloy
B37. Mild Steel	B38. RHA Steel
B39. S7tool steel alloy	B40. AM30 Mg alloy
B41. AM50 Mg alloy	B42. AM60 Mg alloy
B43. AZ31 Mg alloy	B44. AZ91 Mg alloy
B45. AE42 Mg alloy	B46. AE44 Mg alloy
B47. Ti7Al4Mo alloy	B48. Ti8Al1Mo1V alloy
B49. Ti0Al6V4 alloy	B50. Ti6Al6V2Sn alloy
B51. D38 Uranium	B52. D380075Ti Uranium
B53. D38006 Nb Uranium	B54. 200 Nickel
B55. In718 alloy	B56. Copper
B57. OFHC Copper	B58. ETP Copper
B59. 99% Brass	B60. Lead
B61. Tantalum	



B1. Al A356-T6 alloy: model correlation with different temperatures and strain rates

Reference: Horstemeyer, M.F., Lathrop, J., Gokhale, A.M., and Dighe, M., "Modeling Stress State Dependent Damage Evolution in a Cast Al-Si-Mg Aluminum Alloy," *Theoretical and Applied Fracture Mechanics*, Vol. 33, pp. 31-47, 2000.



B2. Al 99% pure alloy: model correlation with different strain rates.

References:

1. Holt, D.L., Babcock, S.G. Green, S.J. and Maiden, C.J., The Strain Rate Dependence of the Flow Stress in Some Aluminum Alloys, General Motors Research Laboratories, TR66-75, 1996.



B3. Al 1060-T0 alloy: model correlation with different strain rates

References:

1. Holt, D.L., Babcock, S.G. Green, S.J. and Maiden, C.J., The Strain Rate Dependence of the Flow Stress in Some Aluminum Alloys, General Motors Research Laboratories, TR66-75, 1996.



B4. Al 1100 alloy: model correlation with different strain rates

References:

1. Holt, D.L., Babcock, S.G. Green, S.J. and Maiden, C.J., The Strain Rate Dependence of the Flow Stress in Some Aluminum Alloys, General Motors Research Laboratories, TR66-75, 1996.

2. Horstemeyer, M.F. and Ramaswamy, S., "On Factors Affecting Localization and Void Growth in Ductile Metals: A Parametric Study," *Int. J. Damage Mech.*, Vol. 9, pp. 6-28, 2000.

3. Johnson, G.R. and Holmquist, T.J., Test data and computational strength and fracture model constants for 23 materials subjected to large strains, high strain rates, and high temperatures, LA-11463-MS, Los Alamos National Laboratory, 1989.



B5. Al 2024-T0 alloy: model correlation with different strain rates

References:

1. Holt, D.L., Babcock, S.G. Green, S.J. and Maiden, C.J., The Strain Rate Dependence of the Flow Stress in Some Aluminum Alloys, General Motors Research Laboratories, TR66-75, 1996.



B6. Al 2024-T351 alloy: model correlation with different strain rates

Reference: Johnson, G.R. and Holmquist, T.J., Test data and computational strength and fracture model constants for 23 materials subjected to large strains, high strain rates, and high temperatures, LA-11463-MS, Los Alamos National Laboratory, 1989.



B7. Al 2024-T4 alloy: model correlation with different strain rates



B8. Al 5083 alloy: model correlation with different strain rates

Reference: Tucker, M.T., Horstemeyer, M.F., Whittington, W.R., Solanki, K.N., Gary, G., and Gullett, P.M., The Effect of Varying Strain Rates and Stress States on the Plasticity, Damage, and Fracture of Aluminum Alloys, Mechanics of Materials, (submitted).



B9. Al 5086-T6 alloy model correlation

Reference: Balasundaram, A., Shan, Z., Gokhale, A.M., Graham, S., and Horstemeyer, M.F., "Particle Rotations During Plastic Deformation of 5086 (O) Aluminum Alloy", *Materials Characterization*, Vol. 48 (no. 5), PP. 363-369, 2002.



B10. Al 6022 alloy model correlation.

Reference: Burton, C., Horstemeyer, M., Wang, P., and Shahbazian-Yassar, R., Rolling history microstructural analysis of texture gradients for 6022 aluminum sheet, *136th TMS Annual Meeting*, Orlando, Fl, Feb. 2007.


B11. Al 6061-T0 alloy: model correlation with different strain rates

1. Holt, D.L., Babcock, S.G. Green, S.J. and Maiden, C.J., The Strain Rate Dependence of the Flow Stress in Some Aluminum Alloys, General Motors Research Laboratories, TR66-75, 1996.



B12. Al 6061-T6 alloy: model correlation with different temperatures and strain rates

1. Department of Defense. Aerospace Structural Metals Handbook, Volume 2. West Lafayette: CINDAS/Purdue, 1993.

2. Holt, D.L., Babcock, S.G. Green, S.J. and Maiden, C.J., The Strain Rate Dependence of the Flow Stress in Some Aluminum Alloys, General Motors Research Laboratories, TR66-75, 1996.1. Horstemeyer, M.F., "A Numerical Parametric Investigation of Localization and Forming Limits," Int. J. Damage Mech, Vol. 9, pp. 255-285, 2000.

3. Johnson, G.R. and Holmquist, T.J., Test data and computational strength and fracture model constants for 23 materials subjected to large strains, high strain rates, and high temperatures, LA-11463-MS, Los Alamos National Laboratory, 1989.



B13. Al 6061-T651 alloy: model correlation with different strain rates, damage state variable

References :

1. Horstemeyer, M.F., "A Numerical Parametric Investigation of Localization and Forming Limits," Int. J. Damage Mech, Vol. 9, pp. 255-285, 2000.



B14. Al 7039 alloy: model correlation with different strain rates



B15. Al 7075-T0 alloy: model correlation with different strain rates

1. Holt, D.L., Babcock, S.G. Green, S.J. and Maiden, C.J., The Strain Rate Dependence of the Flow Stress in Some Aluminum Alloys, General Motors Research Laboratories, TR66-75, 1996.



B16. Al 7075-T6 alloy: model correlation with different strain rates

1. Holt, D.L., Babcock, S.G. Green, S.J. and Maiden, C.J., The Strain Rate Dependence of the Flow Stress in Some Aluminum Alloys, General Motors Research Laboratories, TR66-75, 1996.



B17-a. Al 7075-T651 alloy: damage model correlation

Reference: Jordon, J.B., Horstemeyer, M.F., Solanki, K., Xue, Y., "Damage and stress state influence on the Bauschinger effect in aluminum alloys", *Mechanics of Material*, Vol. 39, pp. 920-931, 2007.



B17-b. Al 7075-T651 alloy: model correlation showing the Bauschinger effect

Reference: Jordon, J.B., Horstemeyer, M.F., Solanki, K., Xue, Y., "Damage and stress state influence on the Bauschinger effect in aluminum alloys", *Mechanics of Material*, Vol. 39, pp. 920-931, 2007.



B18. 300 Maraging steel alloy: model correlation with two strain rates



B19. 1006 steel alloy: model correlation at different strain rates



B20. 1020 steel alloy: model correlation with different strain rates.

1. Nicholas, T., Material behavior at high strain rates, Report AFWAL-TR-80-4053, USAF Wright Aeronautical Laboratories, Wright-Patterson Air Force Base, OH, USA, 1980.

2. Bless, UDRI, 1984.



B21. 10b22 steel alloy: model correlation with different strain rates.



B22. 304L SS alloy: model correlation with different temperatures and strain rates

References:

1. Department of Defense. Aerospace Structural Metals Handbook, Volume 1. West Lafayette: CINDAS/Purdue, 1993.

2. Horstemeyer, M.F., Matalanis, M.M., Sieber, A.M., and Botos, M.L., "Micromechanical Finite Element Calculations of Temperature and Void Configuration Effects on Void Growth and Coalescence," Int J. Plasticity, Vol. 16, 2000.

3. Lu, W. Y., Horstemeyer, M. F., Korellis, J., Grishibar, R., and Mosher, D., "High Temperature Effects in 304L Stainless Steel Notch Tests," *Theoretical and Applied Fracture Mechanics*, Vol. 30, pp. 139-152, 1998.



B23. 321 SS alloy: model correlation with different strain rates.



B24. 410 SS alloy: model correlation with different strain rates



B25. 4140 commercial steel alloy model correlation with different strain rates and temperatures

Reference: Gomez, R.A., Monotonic Plasticity-Damage and Fatigue Life Model Correlations on AISI 4140 Steel, MSc. Thesis, Dept. of Mechanical Engineering, Mississippi State University, 2007.



B26. 4340 steel alloy: model correlation with different strain rates

1. Johnson, G.R. and Holmquist, T.J., Test data and computational strength and fracture model constants for 23 materials subjected to large strains, high strain rates, and high temperatures, LA-11463-MS, Los Alamos National Laboratory, 1989.



B27. A286 steel alloy: model correlation with different strain rates



B28. AF steel alloy: model correlation with different strain rates



B29. Armco Iron : model correlation with different temperatures and strain rates



B30. Armor Steel: model correlation with different temperatures



B31. C1008 Steel : model correlation with different strain rates

1. Nicholas, T., Material behavior at high strain rates, Report AFWAL-TR-80-4053, USAF Wright Aeronautical Laboratories, Wright-Patterson Air Force Base, OH, USA, 1980.

2. Bless, UDRI, 1984.



B32. Carpeleciron: model correlation with different strain rates



B33. FC0205 steel alloy: model correlation with different temperatures

Reference: Allison, P. G., et al., Microstructure-property relations of a powder metallurgy steel (FC-0205) under various loading conditions, (In preparation), 2009.



B34. HY80 steel alloy: model correlation with different temperatures and strain rates



B35. HY100 steel alloy: model correlation with different temperatures and strain rates

References:

1. Johnson, G.R. and Holmquist, T.J., Test data and computational strength and fracture model constants for 23 materials subjected to large strains, high strain rates, and high temperatures, LA-11463-MS, Los Alamos National Laboratory, 1989.

2. Nicholas, T., Material behavior at high strain rates, Report AFWAL-TR-80-4053, USAF Wright Aeronautical Laboratories, Wright-Patterson Air Force Base, OH, USA, 1980.

3. Bless, UDRI, 1984.



B36. HY130 steel alloy: temperature model correlation



B37. Mild Steel: strain rate model correlation

1. Nicholas, T., Material behavior at high strain rates, Report AFWAL-TR-80-4053, USAF Wright Aeronautical Laboratories, Wright-Patterson Air Force Base, OH, USA, 1980.

2. M.J. Manjoine.



B38. RHA Steel: strain rate model correlation



B39. S7tool steel alloy: strain rate model correlation



B40. AM30 Mg alloy model correlation

Reference: M.F. Horstemeyer, D. Oglesby, J. Fan, P.M. Gullett, H. El Kadiri, Y. Xue, C. Burton, K. Gall, B. Jelinek, M.K. Jones, S. G. Kim, E.B. Marin, D.L. McDowell, A. Oppedal, N. Yang, "From Atoms to Autos: Designing a Mg Alloy Corvette Cradle by Employing Hierarchical Multiscale Microstructure-Property Models for Monotonic and Cyclic Loads," MSU.CAVS.CMD.2007-R0001.



B41. AM50 Mg alloy model correlation

Reference: M.F. Horstemeyer, D. Oglesby, J. Fan, P.M. Gullett, H. El Kadiri, Y. Xue, C. Burton, K. Gall, B. Jelinek, M.K. Jones, S. G. Kim, E.B. Marin, D.L. McDowell, A. Oppedal, N. Yang, "From Atoms to Autos: Designing a Mg Alloy Corvette Cradle by Employing Hierarchical Multiscale Microstructure-Property Models for Monotonic and Cyclic Loads," MSU.CAVS.CMD.2007-R0001



B42. AM60 Mg alloy: temperature model correlation

Reference: M.F. Horstemeyer, D. Oglesby, J. Fan, P.M. Gullett, H. El Kadiri, Y. Xue, C. Burton, K. Gall, B. Jelinek, M.K. Jones, S. G. Kim, E.B. Marin, D.L. McDowell, A. Oppedal, N. Yang, "From Atoms to Autos: Designing a Mg Alloy Corvette Cradle by Employing Hierarchical Multiscale Microstructure-Property Models for Monotonic and Cyclic Loads," MSU.CAVS.CMD.2007-R0001



B43. AZ31 Mg alloy: temperature and strain rate model correlation

Reference: F.J. Polesak, III, C.E. Dreyer, T.J. Shultz, and S.R. Agnew, "Blind Study of the Effect of Processing History on the Constitutive Behavior of Alloy AZ31B," Magnesium Technology 2009, Eds. E.A. Nyberg, S.R. Agnew, M.O. Pekguleryuz, and N.R. Neelameggham (TMS-AIME: Warrendale, PA: 2009) 491-496.



B44. AZ91 Mg alloy: temperature model correlation

Reference: M.F. Horstemeyer, D. Oglesby, J. Fan, P.M. Gullett, H. El Kadiri, Y. Xue, C. Burton, K. Gall, B. Jelinek, M.K. Jones, S. G. Kim, E.B. Marin, D.L. McDowell, A. Oppedal, N. Yang, "From Atoms to Autos: Designing a Mg Alloy Corvette Cradle by Employing Hierarchical Multiscale Microstructure-Property Models for Monotonic and Cyclic Loads," MSU.CAVS.CMD.2007-R0001



B45. AE42 Mg alloy: temperature and strain rate model correlation

Reference: M.F. Horstemeyer, D. Oglesby, J. Fan, P.M. Gullett, H. El Kadiri, Y. Xue, C. Burton, K. Gall, B. Jelinek, M.K. Jones, S. G. Kim, E.B. Marin, D.L. McDowell, A. Oppedal, N. Yang, "From Atoms to Autos: Designing a Mg Alloy Corvette Cradle by Employing Hierarchical Multiscale Microstructure-Property Models for Monotonic and Cyclic Loads," MSU.CAVS.CMD.2007-R0001


B46. AE44 Mg alloy: temperature model correlation

Reference: M.F. Horstemeyer, D. Oglesby, J. Fan, P.M. Gullett, H. El Kadiri, Y. Xue, C. Burton, K. Gall, B. Jelinek, M.K. Jones, S. G. Kim, E.B. Marin, D.L. McDowell, A. Oppedal, N. Yang, "From Atoms to Autos: Designing a Mg Alloy Corvette Cradle by Employing Hierarchical Multiscale Microstructure-Property Models for Monotonic and Cyclic Loads," MSU.CAVS.CMD.2007-R0001



B47. Ti7Al4Mo alloy: strain rate model correlation



B48. Ti8Al1Mo1V alloy: strain rate model correlation



B49. Ti0Al6V4 alloy: strain rate model correlation

References:

1. Department of Defense. Military Standardization Handbook: Metallic Materials and Elements for Aerospace Vehicle Structures MIL-HDBK-5E, Volume 2. June 1987.

2. Johnson, G.R. and Holmquist, T.J., Test data and computational strength and fracture model constants for 23 materials subjected to large strains, high strain rates, and high temperatures, LA-11463-MS, Los Alamos National Laboratory, 1989.

3. Nicholas, T., Material behavior at high strain rates, Report AFWAL-TR-80-4053, USAF Wright Aeronautical Laboratories, Wright-Patterson Air Force Base, OH, USA, 1980.

4. Maiden and Green



B50. Ti6Al6V2Sn alloy: strain rate model correlation

References:

1. Department of Defense. Military Standardization Handbook: Metallic Materials and Elements for Aerospace Vehicle Structures MIL-HDBK-5E, Volume 2. June 1987.



B51. D38 Uranium: strain rate model correlation



B52. D380075Ti Uranium: strain rate model correlation



B53. D38006 Nb Uranium: strain rate model correlation



B54. 200 Nickel: temperature model correlation



B55. In718 alloy: strain rate model correlation



B56. Copper: strain rate and temperature model correlation.

Reference: Tanner A.B., McGinty R.D. and McDowell D.L., Modeling temperature and strain rate history effects in OFHC Cu, Int. Journal of Plasticity, 1999.



B57. OFHC Copper: strain rate model correlation

References:

1. Johnson, G.R. and Holmquist, T.J., Test data and computational strength and fracture model constants for 23 materials subjected to large strains, high strain rates, and high temperatures, LA-11463-MS, Los Alamos National Laboratory, 1989.

2. Nicholas, T., Material behavior at high strain rates, Report AFWAL-TR-80-4053, USAF Wright Aeronautical Laboratories, Wright-Patterson Air Force Base, OH, USA, 1980.

3. U.S. Lindholm



B58. ETP Copper: strain rate model correlation



B59. 99% Brass: temperature and strain rate model correlation



B60. Lead: strain rate model correlation

References:

1. Nicholas, T., Material behavior at high strain rates, Report AFWAL-TR-80-4053, USAF Wright Aeronautical Laboratories, Wright-Patterson Air Force Base, OH, USA, 1980.

2. U.S. Lindholm



B61. Tantalum: temperature and strain rate model correlation