Quarterly Management Document FY20, 2nd Quarter, Physics-based Creep Simulations of Thick Section Welds in High Temperature and Pressure Applications

Thomas M Lillo, Wen Jiang

April 2020



The INL is a U.S. Department of Energy National Laboratory operated by Battelle Energy Alliance

Quarterly Management Document FY20, 2nd Quarter, Physics-based Creep Simulations of Thick Section Welds in High Temperature and Pressure Applications

Thomas M Lillo, Wen Jiang

April 2020

Idaho National Laboratory Idaho Falls, Idaho 83415

http://www.inl.gov

Prepared for the
U.S. Department of Energy
Office of Fossil Energy
Under DOE Idaho Operations Office
Contract DE-AC07-05ID14517

Quarterly Management Document – FY20, 2nd Quarter, Physics-based Creep Simulations of Thick Section Welds in High Temperature and Pressure Applications

Document # INL/EXT-20-58120

WBS Element	Project Title	Contract Number		Contract Start	Contract End
C.B.10.02.02.4	Physics-based Creep Simulations of	FEAA90		10/01/17	09/30/2020
0	Thick Section Welds in High				
	Temperature and Pressure				
	Applications				
Performer Name and A		Principal Investigator(s)			
Thomas Lillo		Tho	mas Lillo		
Idaho National La					
P.O. Box 1625					
Idaho Falls, ID 83	415				

BUDGET AND COST REPORT

Prior Year Funds	s (\$K)			10.9								
Total Current Ye	ar Commi	tment (\$K)		10.9								
Projected Curre	nt Year Co	sts (\$K)		10.9								
	0	N	D	J	F	М	Α	М	J	J	Α	S
Monthly Planned Costs	0.5	1.0	1.0	1.0	2.7	2.7	2.0	0	0	0	0	0
Actual Monthly Costs	0.6	0.1	1.1	1.1	0.0	0.2						
Monthly Variance	-0.1	0.9	-0.1	-0.4	-2.7	-2.5						
Total costs – planned	0.5	1.5	2.5	3.5	6.2	8.9	10.9	10.9	10.9	10.9	10.9	10.9
Total costs - actual	0.6	0.7	1.8	2.9	2.9	3.1						

MILESTONE REPORT

Milestone Designation	Milestone Description	Due Date	Revised Due Date	Completion Date
A	Evaluate current MOOSE capabilities	09/30/2015		09/30/2015
В	Complete Alloy 617 weld characterization	10/30/2015		11/18/2015
С	Receipt of Alloy 740H plates	10/30/2015		11/05/2015
D	Complete welds in Alloy 740H	11/16/2015	7/31/2016	7/31/2016

Е	Characterize Alloy 740H welds	02/01/2016	09/30/2016	9/02/2016
F	Creep model development – Stage 1	09/30/2016		9/30/2016
G	Creep Model Development – Stage 2	8/29/2017	2/28/2019	1/15/2019
Н	Calibration of Secondary creep – Alloy 617	9/30/2017	3/31/2019	Eliminated
Ι	Stress Drop Tests	2/01/2017	5/31/2018	6/28/2018
J	Characterization of creep failure mechanisms	4/01/2017	04/30/2018	5/04/2018
K	Secondary creep calibration for welds – Alloy 617	5/30/2018	4/15/2019	Eliminated
L	Creep model development – Completion of Stage 3	8/30/2018	8/16/2019	9/9/2019
M	Creep simulation of a welded joint in Alloy 740H	9/30/2018	08/31/2020	
N	Validation of creep simulation model via an Alloy 740H weld consisting of refined microstructure	9/15/2018	09/30/2020	

TECHNICAL HIGHLIGHTS

Milestone M, "Creep simulation of a welded joint in Alloy 740H"

Calibration of the model continued using the experimental creep data for the base metal. A few errors in the execution of the model were discovered and corrected. The model was then calibrated using base metal creep data from the highest temperature (800°C) and the lowest temperature (700°C) creep data with the intent of comparing model simulations to creep data from 750°C. Figures 1 and 2 show the model simulations for 700°C, 413 MPa and 800°C, 200 MPa, respectively.

Most notably in Fig. 1, the model simulates the primary and secondary creep behavior quite well but deviates considerably in the tertiary regime – tending to overestimate the creep strain and, thus, underestimating the creep rupture life by about a factor of 2 (however, the simulation was not run to failure – it is estimated that the simulated creep life will be significantly shorter than the experimental creep life). In the world of creep data, this is not unexpected due to variations in microstructure and experimental testing complexities. However, since the model was calibrated using this experimental data, it would intuitively seem that a better fit to the experimental data should be obtained. The tertiary regime has been an issue since the physical mechanism responsible for the transition to tertiary creep has not been definitively identified (Lillo, T.M., Wright, R.N., "The Onset of Tertiary Creep in Alloy 617", proceedings of the 2015 PVP Pressure Vessel and Piping Conference, Boston, MA, July 19-23, 2015.). We have chosen to use a damage model described by Shen (C. Shen, Modeling Long-term Creep Performance for Welded Nickel-base Superalloy Structures for Power Generation Systems, DOE/NETL Cooperative Agreement DE-FE0024027, 2015) and consists of damage due to dislocation motion as well as damage due to diffusional processes (see, the 3rd Quarter Management Report, FY19 for further details). The damage parameter for the 700°C, 413 MPa simulation also has been plotted in Fig. 1 (the gold line), and, again, in Fig. 3 by itself. The damage parameter is largely linear throughout about

half of the creep life and then starts to accelerate toward the end of life (i.e., deviates from linearity) which is intuitively expected. However, at high values of the damage parameter, i.e., toward the end of life, the rate of damage accumulation in the simulation *decreases*. This is not expected as the damage increases the effective stress (damage equates to loss of load-bearing cross-sectional area in the Shen damage model) in the gage section of the creep specimen and thus the damage rate is expected to accelerate as the specimen reaches end-of-life. The result of the decreasing damage rate is manifested as a decrease in the creep rate at the end of life as shown in Fig. 1. It is not clear at this time the cause behind this behavior but it is thought this is not correct and efforts are underway to understand and, if necessary, correct this behavior in the model.

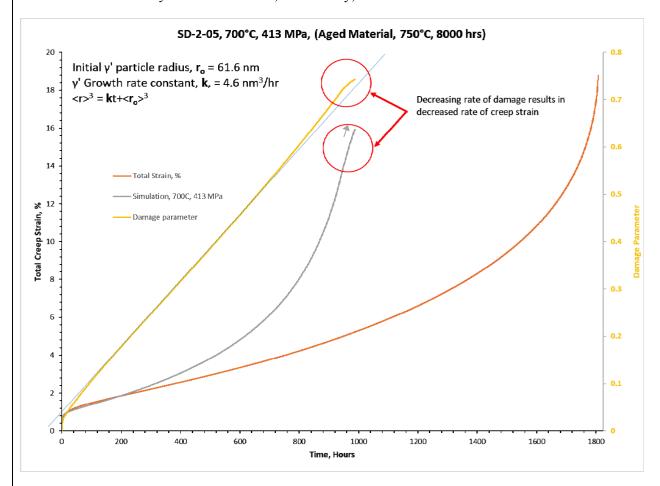


Figure 1. Plots of the experimental creep behavior of the base metal at 700°C (orange curve) with the simulation from the model overlaid (grey curve). The damage parameter from the simulation is also plotted (gold curve). The base metal material for sample SD-2-05 was aged for 8000 hrs at 750°C prior to fabrication of the creep specimen.

At 800°C, the model, again, predicts primary and secondary creep quite well, Fig. 2. In this case, however, the simulation significantly underestimates the tertiary creep strain. The simulation was not run to high values of the damage parameter, which indicates end-of-life, and the damage parameter exhibits linear behavior over the time of the simulation. The damage parameter in this simulation does not exceed about 0.37 – compared to about 0.74 for the simulation in Fig. 1.

After calibration using the experimental data in Figs. 1 and 2, the model was used to simulate the creep behavior at 750°C and 305 MPa, Fig. 4. The model predicts primary and secondary creep behavior under these

conditions quite well when compared to the two experimental datasets at these conditions. (The experimental data does show considerable variation in creep behavior with the creep life varying by about 25% or a factor of 1.25.) However, again, the creep strain associated with tertiary creep is underestimated. Again, the simulation was not run to end-of-life (i.e., high damage values) but it would appear the model predicts a value for end-of-life that is approximately a factor of 2 greater than the actual creep life of the experimental creep tests. The damage parameter shows deviation from linearity in Fig. 4 (gold curve), but, again, the simulation was only run to a maximum damage parameter value of about 0.55 (in theory, a damage parameter of 1 indicates failure and end-of-life).

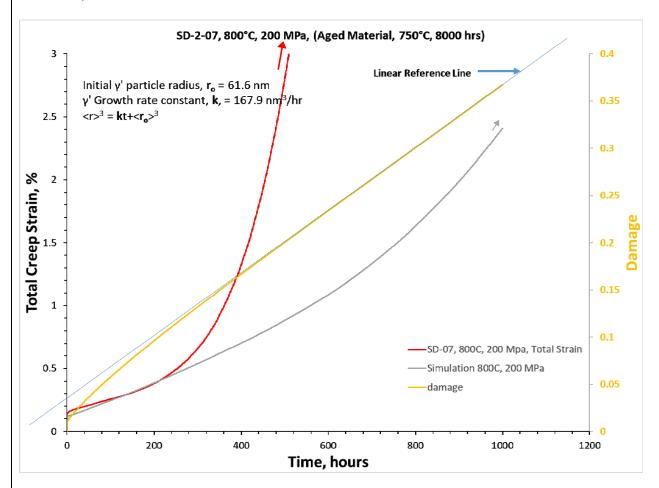


Figure 2. Plot of the experimental creep behavior at 800°C and 200 MPa for base metal (red curve) and the simulated creep behavior (grey curve). The damage parameter is also plotted in this figure (gold curve). The material for the base metal sample, SD-2-07, was aged for 8000 hrs at 750°C prior to fabrication of the creep specimen.

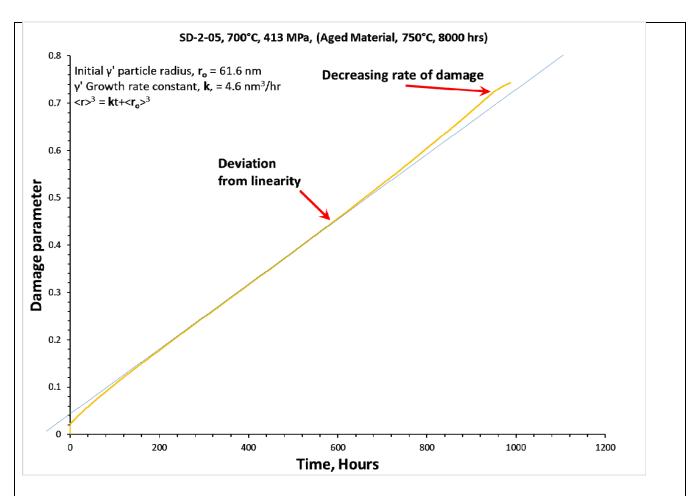


Figure 3. Plot of the damage parameter for the simulation of 700°C and 413 MPa. The damage parameter deviates from linearity later in the simulated creep life, as expected, but also exhibits a decreasing damage rate at the very end of life which is not expected.

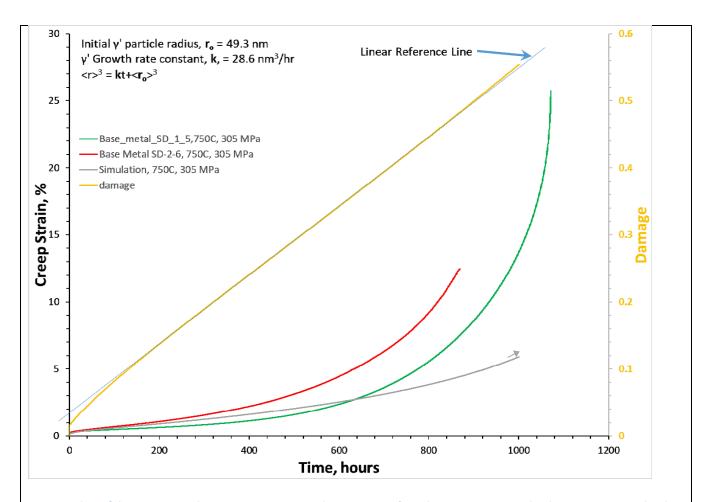


Figure 4. Plots of the experimental creep curves, SD-1-5 and SD-2-6, at 750°C and 305 MPa, green and red curves, respectively. The simulated creep curve (grey curve) is also plotted. Finally, the damage parameter is included in the figure.

Overall, the model simulates primary and secondary creep under the various experimental creep conditions. Figures 1, 2 and 4 seem to indicate that damage is overestimated at 700°C, 413 MPa and underestimated at the 750 and 800°C creep conditions. The influence of diffusion increases with test temperature, which, experimentally, ranges from about 0.64 T_{MP} to about 0.70 T_{MP}, and can be expected to influence the contribution of diffusion to the simulation. Since the model simulates primary and secondary creep well for each test temperature, it is likely the influence of diffusion on the damage parameter evolution is not quite as it should be. A closer look at the influence of diffusion on the evolution of the damage parameter will be carried out early next quarter. Additionally, changes to the calibration methodology will be explored, e.g., calibration using the two sets of creep data taken at 800 and 750°C versus the current calibration using the 800 and 700°C data.

Completion of this task is now targeted for 08/31/2020.

Milestone N, "Validation of creep simulation model via an Alloy 740H weld consisting of refined microstructure"

No progress on this task during the 2nd quarter of FY20.

The completion date of this task is now targeted for 09/30/2020

ISSUES

Tertiary creep remains a hurdle to be overcome in the simulation. The reason for this is due the ambiguous mechanism resulting in the transition to tertiary creep. It is likely a physical mechanism is responsible for the transition to tertiary creep and it remains to be seen whether the damage parameter evolution approach will prove a satisfactory description of the mechanism(s) responsible for the transition to tertiary creep. At this point, we will look at this to the extent possible (i.e., within the limits of the remaining funds) and then move on to simulating the all-weld metal creep tests. After this, the simulation of cross weld creep tests will be a combination of the base metal behavior and the weld metal behavior based on a rule-of-mixtures approach based on the fractions of base metal and weld metal in the gage section of the creep specimen.

Report Prepared By	Date
Thomas M. Lillo and Wen Jiang	04/28/2020