

## Article

# Application of mathematical search algorithms for unknown material properties in Additive Manufacturing simulations

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**Abstract:** Additive manufacturing (AM) simulations have risen as a way to better understand the effect of processing parameters on builds. They are effective for materials which are well characterized

and published, however for newer or proprietary materials, they cannot provide accurate results due to the lack of knowledge of material properties. This work demonstrates the process of the application of mathematical search algorithms to develop an optimized material dataset which results

in accurate simulations. This was done with 7000 series aluminum and the laser directed energy deposition (DED) process. The Nelder-Mead search algorithm was able to develop an optimized dataset which had a combined width and depth error of just 9.1%. This optimization started from a generic aluminum material properties dataset found in the literature which had an initial error of

600%. Demonstrating that it is possible to develop an optimized material property dataset which facilitates a more accurate simulation for an undercharacterized material.

**Keywords:** Additive manufacturing (AM), Mathematical modeling, Mathematical search, Material properties, Aluminum, Additive manufacturing (AM) simulation, Input parameter optimization

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## 1. Introduction

Additive manufacturing (AM) is an emerging manufacturing technique which has the potential to revolutionize manufacturing. To realize this revolution, it is necessary to be able to produce components reliably and to understand the process well enough to ensure that builds are consistent enough that the performance of the completed build can

<sup>1</sup> Updated the keywords

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be guaranteed. To do this, researchers and manufacturers have turned to mathematical modeling to understand the process [1].

The differences in simulation techniques can vary based on the desired response from the simulation and the underlying assumptions which were made during the development of the models. One example of this can be seen when comparing the mathematical models presented in [1], [2], and [3]. They all attempt to model roughly the same aspect of the build but take very different approaches. [1] take a purely physics-based approach to the solution and works from first principle of the physical process being modeled. On the contrary, [2] is a data driven model which used a breadth of data to develop a mathematical model which properly predicts the material behavior. In between these models exists [3] which attempts to marry the two approaches and develop a physics-based model that uses data to improve accuracy. The one unifying characteristic of these, and all, mathematical models, is the need for the inclusion of a dataset which defines the behavior of the material being investigated, this is colloquially referred to as the material properties. These material properties can vary in literature and this variance in values can lead to a discrepancy in simulation results [?].

Though variation exists in the literature, values can be found and used when the material is well characterized and published. An example of a well published material is Ti-64 where it is easy to find literature which reports the material properties such as in [4], [5], and [6]. However, for materials which are not well understood and published, such as specific aluminum alloys, it can be very challenging though possible to find some properties such as [7] and [8]. There is a need to determine a complete dataset of the material properties, or at a minimum, develop the dataset which produces the most accurate simulation results. This can be accomplished by expending the necessary resources to measure the needed properties using advanced equipment. This process can be expensive and time-consuming which has led to the development of material simulations which attempt to predict the material properties, such as [9]. Though faster and cheaper than experimental results, they, like all simulations, have an error that is associated with them. Using these values alone can lead to unknown errors stacking up in the AM models. This problem also applies to materials where the tolerances for alloying elements is so wide that specimen of the same alloy can have different material properties.

In order to develop a dataset which produces accurate AM simulation results, a multi-objective optimization scheme can be used as a search algorithm to determine the dataset *Routine*.

*Accuracy & Material*

*Put for JMatPro?*

*from JMatPro*

More references  
about materials  
Simulator more  
detailed

This work will perform  
on optimization routine  
on the material dataset  
for the 7050 Al alloy to  
enable more accurate simulation  
results. This will be  
done by validating the  
model on the well characterized  
Ti-64 material showing its  
models effectiveness. Then  
utilizing sister alloys,  
a generic material properties  
dataset will be found from  
literature. The  
search algorithm will be applied  
to generate an optimal input  
properties set. These material input  
will then be used  
with a range of laser power  
and confirm experimentally to  
show effectiveness over a  
processing window.

which produces the most realistic results. This can be used to develop a dataset for a new alloy along with for a specific batch of stock which has been procured from a supplier.

This work will aim to address one of the current desires in metal AM which is to be able to produce parts out of aluminum. The desire is evident by the volume of effort being applied to aluminum AM ([10], [11], [12]). The challenge associated with this stems from the wide range of alloys which have wildly varying material properties and are not well published. One of the weldable high-strength alloys which has been targeted for metal AM is 7050 [13]. Though the material is widely available, temperature-dependent material properties are not readily available. Therefore, this work will find material properties in literature which are an approximation of the 7050 aluminum, namely sister alloys with similar compositions, as a starting point for the search algorithm. From that starting database, a search algorithm will be applied to develop a material dataset which produces more accurate AM simulation results.

## 2. Materials and Methods

### RESUME

This body of work begins by describing the thermal model which was developed to simulate the metal AM deposition process for any metal where the material properties could be determined. This model is then validated using a well characterized material, Ti-64. Upon validation of the thermal model, it is applied to a generic aluminum alloy and experimentation is used to determine if the results are a good characterization of the physical process. When found to be inaccurate, a mathematical search algorithm was applied to the input parameters in order to increase the accuracy of the models results. Lastly, the optimal input dataset was validated by comparing the simulated results to experimental results with a range of laser scan speeds and power levels.<sup>2</sup>

### 2.1. Simulation description

The model used in this study was developed at Missouri University of Science and Technology. This simulation has the express goal of being efficient while still holding true to physics models. In order to accomplish this, it heavily leverages GPU processing by utilizing image processing techniques. The simulation was developed with the laser DED processes in mind, however, it was developed in a modular manner such that it can be applied to most AM processes.

The model is a voxel-based simulation, which forgoes the calculation of the fluid flow and focuses on heat transfer and material insertion. This decision was based on past

<sup>2</sup> This section was added to provide a sketch of the methods used in this work

simulation development experience, where it was determined that the calculation of the fluid flow was the most computationally expensive part of the simulation.

The governing equations of the model developed can be seen in Equations 1, 2, and 3 which describe the flow of heat from conduction, convection and radiation, and laser absorption respectively [14].

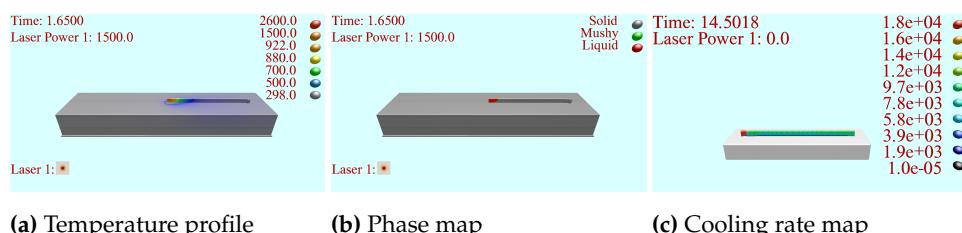
$$\rho c_p \frac{\partial T}{\partial t} = k \left( \frac{\partial^2 T}{\partial x^2} + \frac{\partial^2 T}{\partial y^2} + \frac{\partial^2 T}{\partial z^2} \right) \quad (1)$$

$$k \frac{\partial T}{\partial n} = -h(T_s - T_a) - \varepsilon \sigma (T_s^4 - T_a^4) + \phi(x, y, z) + \rho c \frac{\partial T}{\partial t} \quad (2)$$

$$\phi(x, y, z) = H(z) \alpha \phi_0 \sqrt{1 - \frac{x^2}{r_0^2} - \frac{y^2}{r_0^2}} \quad (3)$$

In these equations,  $\rho$  is density,  $c_p$  is specific heat,  $k$  is thermal conductivity,  $h$  is convection coefficient,  $\varepsilon$  is emissivity,  $\sigma$  is Stefan-Boltzmann constant,  $n$  is the unit (outward) normal vector of a point at location  $(x, y, z)$  that is located on the outer surface of the component,  $r_0$  is the radius of the laser beam,  $\alpha$  is the absorption of the material with respect to the laser radiation,  $T_s$  is the surface temperature,  $T_a$  is the ambient temperature,  $\phi_0$  is the laser power, and  $H(z)$  is a step function which is 1 for the node with the largest  $z$  value in every  $(x, y)$  location and 0 elsewhere.

The main attributes of the model are its ability to predict the thermal history of a part, Figure 1a, the phase map of the part at any given time, Figure 1b, and a cooling rate for any section of the part, Figure 1c. This helps to give a predictor of the microstructure within the deposition which is the driving force behind the final mechanical properties. Additionally, the laser in the simulation is modeled in 3-D to be able to take into account



**Figure 1.** Examples of data maps which can be expected from the simulation.

the beam quality using the beam parameter product (BPP) reported by the manufacturer and shown in Equation 4, where  $\theta$  and  $w_0$  are the divergence angle and the beam waist respectively.

$$BPP = 0.5\theta w_0 \quad (4)$$

Lastly, the model includes true ray tracing enabling shadowing of the laser, the ability to define a mass which represents the machine acting as a heat sink during the build process, and the inclusion of temperature-dependent material properties.

The main objective of the model is to accurately predict the thermal history of the build. In order to initially validate the models, the well-characterized and published material of Ti-64 was used. The validation was done by scanning a laser on the surface of a substrate at three energy densities with the experimental parameters found in Table 1 and the material properties used in the simulations can be seen in Table 2.<sup>3</sup>

**Table 1.** Simulation parameters used in Ti-64 validation

Parameter	Value
Resolution	60 $\mu\text{m}$
Laser diameter	2.0 mm
Laser Profile	TEM00
Laser power	1000 W
Energy density (Equation 5)	13, 18, 24 $\frac{\text{W}}{\text{mm}^2/\text{sec}}$
Scan Length	45 mm
Substrate dimensions	55 mm x 12.7 mm x 6.35 mm

**Table 2.** Ti-64 material properties used in validation

Material Property	Value	Reference
Solidus temperature	1603°C	[4]
Liquidus temperature	1650°C	[15]
Solid density	4420.0 $\frac{\text{kg}}{\text{m}^3}$	[15]
Fluid density	3920.0 $\frac{\text{kg}}{\text{m}^3}$	[15]
Specific heat	0.4 - 1.15 $\frac{\text{J}}{\text{gK}}$	[5]
Thermal conductivity	5.0 - 43.0 $\frac{\text{W}}{\text{mK}}$	[5]
Absorptivity	0.4	[6]

There are several representations of the energy density of a laser beam in AM but the most appropriate method of calculating the energy for the DED process is to use a surface energy density, Equation 5, where P is laser power, A is the laser spot area, and v is the scan speed [16].

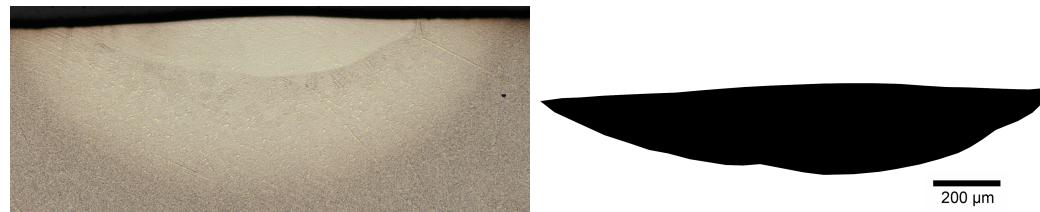
$$SED_{area} = \frac{P}{Av} \quad (5)$$

In order to analyze the results, the samples from the three parameter sets were sectioned using a wire electrical discharge machine (EDM), polished using an automated polisher to a mirror finish, and etched using Kroll's reagent<sup>4</sup> to make the melt track visible with an optical microscope as can be seen in Figure 2a, this was done at three points along

<sup>3</sup> Updated table to include the property values for Ti-64

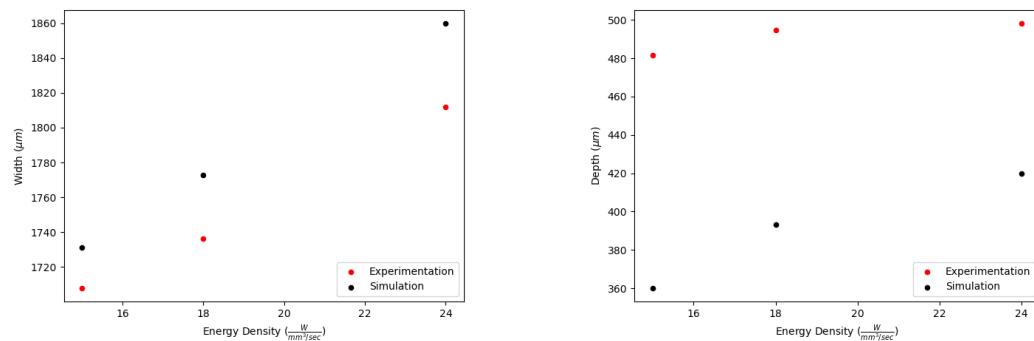
<sup>4</sup> Detailed Ti64 etching routine

the melt length. The experiment and simulation were compared and the resulting error



**Figure 2.** Analysis of sliced sample in Ti-64 validation

for the width can be seen in Figure 3a and the depth can be seen in Figure 3b. The average



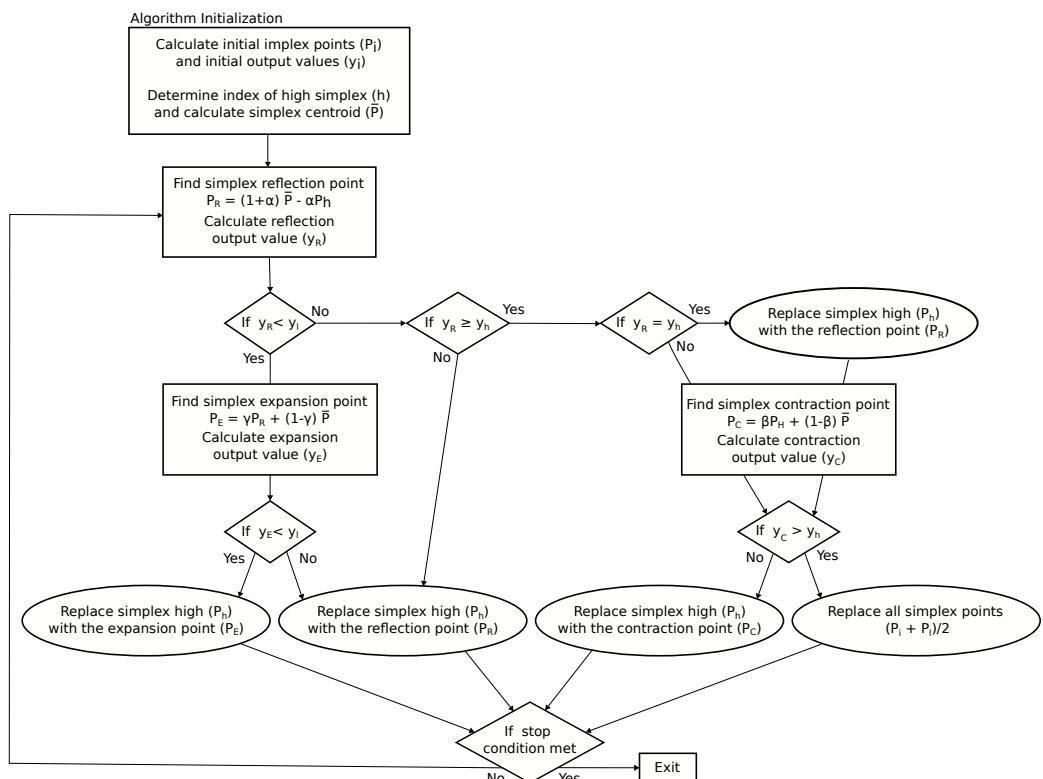
**Figure 3.** Comparison of simulation and experimentation in Ti-64 validation.

values which are reported on the graph have an associated standard deviation of the  $0.62 \mu m$ ,  $13.7 \mu m$ , and  $7.74 \mu m$  for the width measurements for the energy densities of  $15 \frac{W}{mm^3/sec}$ ,  $18 \frac{W}{mm^3/sec}$ , and  $24 \frac{W}{mm^3/sec}$  respectively. For the track depth the standard deviations are  $70.2 \mu m$ ,  $113.7 \mu m$ , and  $13.7 \mu m$  for the energy densities of  $15 \frac{W}{mm^3/sec}$ ,  $18 \frac{W}{mm^3/sec}$ , and  $24 \frac{W}{mm^3/sec}$  respectively.

From the plots, it can be seen that the simulation is capable of predicting the width within 3% error and the depth error is between 15% and 25%. The error in the width is very acceptable at less than 3%. However, the error in the depth is larger due to the resolution (voxel size) of the simulation chosen and its relative size to the depth. This error, of approximately 20%, for the depths which range from  $450-500 \mu m$  is 1.3-1.7 times the resolution,  $60 \mu m$ , of the simulation. Since the trends of the simulation and experimentation match and the error is within two resolution distances the 20% error in the depth is considered acceptable as well. This shows that the mathematical models developed are accurate and if the material property is well characterized, the simulation will produce accurate results.

## 2.2. *SearchTuning* algorithm description

The search algorithm chosen was the Nelder-Mead search algorithm [17]. This method was selected because it is one of the most popular direct search methods for the minimization of functions. The Nelder-Mead approach is a local optimization search which does not rely on knowledge of the gradient to select the next search point. This is critical for the application to simulation results because the gradient is unknown and finding it would involve running a large number of simulations. With simulation times that can reach into days long, this is a critical consideration. Instead of knowing the actual function, it relies on  $n+1$  vertices. This results in a smaller number of simulation runs being needed to perform the minimization [18]. The flow chart in Figure 4 is the flow which is used to determine the next search point.



**Figure 4.** Flow chart for the Nelder-Mead search algorithm

The search begins by calculating the reflection point using Equation 6 where  $\alpha$  is the reflection coefficient.

$$P_{refl} = (1 + \alpha)P_{cent} - \alpha P_{high} \quad (6)$$

If this reflection point is smaller than the smallest current simplex value, then the expansion is calculated using Equation 7, where  $\gamma$  is the expansion coefficient.

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$$P_{exp} = \gamma P_{refl} - (1 - \gamma) P_{center} \quad (7)$$

If the expansion point is smaller than the reflection point, then the expansion point is used to replace the largest simplex member. Otherwise, if the reflection point is larger than the expansion point, the reflection point is used to replace the largest member of the simplex, and the algorithm is restarted. If the reflection point is larger than the smallest simplex point and smaller than the second largest point, then the highest point of the simplex is replaced with the reflection and the algorithm is restarted. If the reflection point is between the simplex highest value and second highest value, a contraction is calculated, using Equation 6, with the highest values being replaced with the reflection. Otherwise the contraction is calculated with the original simplex still using Equation 6, where  $\beta$  is the contraction coefficient.

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$$P_{cont} = \beta P_{high} - (1 - \beta) P_{cent} \quad (8)$$

If the contraction point is smaller than the largest point of the simplex, then the contraction replaces the largest point and the algorithm is continued. However, if the contraction point is larger than the highest point, a shrink step is performed, detailed in Equation 9, where  $\delta$  is the shrink coefficient and the algorithm is restarted.

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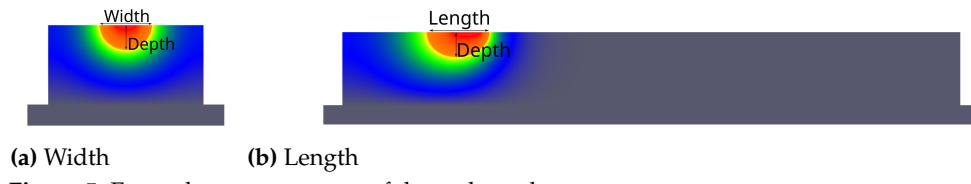
$$P_i = \delta P_i + (1 - \delta) P_{low} \quad (9)$$

### 2.3. Selection of Properties

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One of the attractive characteristics of the Nelder-Mead search algorithm is its ability to scale to an unlimited number of unknowns. The main adverse effects of the scaling are the increased number of runs and the combination of the errors. As the number of unknowns increases the complexity of the search space increases, which in turn increases the number of iterations needed to find a minimum. This can result in drastically longer wait times for the search results. Additionally, with the increased number of unknowns, the stop condition of the search algorithm will be met at a different interval since the stop condition is based on the variance in the simplex. This may result in modifications to the stop conditions being necessary as a larger number of unknowns are included. If the response variable is properly defined, it will not affect the model's results if more unknowns are included [18].

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**Figure 5.** Example measurements of the melt pool

In order to reduce the complexity of the search algorithm, a sensitivity analysis was performed. This work began by finding the material properties which were needed in the models as summarized in Table 3.

**Table 3.** Key material properties in thermal modeling of AM

Material Property	Reference
Solidus temperature	[19], [7], [20], [21]
Liquidus temperature	[19], [7], [20], [21]
Solid density	[21], [22]
Fluid density	[21], [23], [8]
Specific heat	[7], [8]
Thermal conductivity	[7], [8]
Absorptivity	[24], [25], [26]

These properties were varied according to a Plackett-Burman design of experiment in order to determine the properties which, when changed, had a statically significant effect on the resulting melt pool width, depth, and volume, as measured in Figure 5. Analyzing these results with Pareto charts of the standardized effects of the variables and partial regression plots of the residuals it was determined that the variables in Table 4 had a statically significant effect on the resulting melt track when modified. This work included

**Table 4.** Critical material properties

Laser absorption at 880°C
Laser absorption at 922°C
Thermal conductivity at 922°C
Thermal conductivity at 1491°C
Specific heat at 733°C

the laser diameter in the search algorithms dataset due to the difficulty associated with accurately measuring the diameter, this process is detailed in [27].

#### 2.4. Parameter Search Tuning and Simulation setup

For this study, the Nelder-Mead search algorithm parameters which were used can be seen in Table 5. These parameters were chosen because they fell within the guidelines from the algorithm description and after trial and error produced the most efficient searching tuning [17].

**Table 5.** Nelder-Mead algorithm parameters

Parameter	Value
$\alpha$	5.0
$\gamma$	10.0
$\beta$	0.5
$\sigma$	0.5

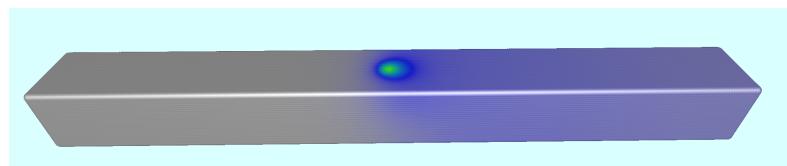
The starting values which were used as the starting point for the search algorithm can be seen in Table 6. These values were chosen based on the values which were found in the literature for similar aluminum alloys. The laser diameter was chosen based on the measuring of a melt track width on a substrate. To ensure the search algorithm did not

**Table 6.** Generic Aluminum materials properties found in the literature

Property	Material Temp.	Value	Ref.
Laser absorption	880°C	15.0%	[25]
Laser absorption	922°C	30.0%	[25]
Thermal conductivity	922°C	88.8 $\frac{W}{mK}$	[8]
Thermal conductivity	1491°C	104.9 $\frac{W}{mK}$	[8]
Specific heat	733°C	1108.0 $\frac{J}{kgK}$	[8]
Laser diameter		1.6 mm	

waste time searching in unacceptable regions the constraint was included that none of the material properties were allowed to become negative. This was the only constraint which was needed in addition to the inherent constraints of the Nelder-Mead algorithm.

The experimental setup which was used was a simple laser scanning of the surface of a substrate, as shown in Figure 6, with the parameters shown in Table 7. This was chosen in order to simplify the experiment. This setup removes the complexity associated with adding material which includes the rate of material addition, molten metal flow parameters, and acceleration effect associated with turning during deposits.

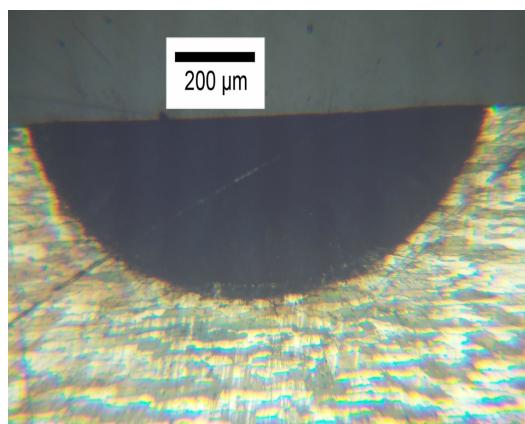
**Figure 6.** Example of simulation setup used to determine melt track size**Table 7.** Experimental constants used in search algorithm tuning experiments

Parameter	Value
Resolution (voxel size)	100 $\mu m$
Laser Power	1,750 W
Laser Scan Speed	1143 mm/min
Laser Profile	Top Hat
Scan Length	77 mm
Substrate dimensions	82 mm x 8 mm x 8 mm

## 2.5. Simulation analysis

Upon completion of each simulation run, the saved data files were analyzed to determine the regions of the simulation which had melted. This was done by developing a map which marked the locations of the domain that had ever been in the fluid phase. This map was then used to determine the width and depth of the melt track along the scan length, excluding the beginning and end where effects from starting and stopping motion would affect the results. These width and depth measurements along the scan length were averaged to develop a single measurement which could be compared to experimentation.

The experimental results were similarly determined, however instead of a continuous set of measurements, there were 4 discrete measurements. These were obtained by slicing the substrate at the prescribed locations using a wire EDM. These slices were polished using an automatic polishing machine up to a mirror finish and etched using Keller's reagent<sup>5</sup> in order to make the microstructural differences visible by an optical microscope, an example of this can be seen in Figure 7, where the dark region had been melted during the experimentation.



**Figure 7.** Example of sliced, polished, and etched slice from Aluminum experimentation

In order for the Nelder-Mead search to function properly, a response variable needed to be defined. This function needed to characterize the accuracy of the simulation into a single parameter which could be minimized and upon minimization would result in the most accurate simulation. To accomplish this goal, Equation 10 was developed. This equation takes into account the error in the width of the simulation along with the error in the depth. This equation results in a non-negative number where 0 represents a simulation which perfectly matched experimentation.

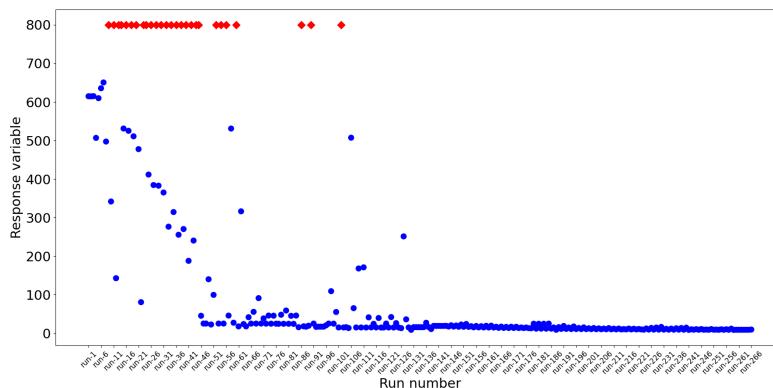
<sup>5</sup> Detailed Al etching routine

$$\text{Response} = \left( \frac{|Sim. Width - Exp. Width|}{Exp. Width} + \frac{|Sim. Depth - Exp. Depth|}{Exp. Depth} \right) * 100 \quad (10)$$

### 3. Results

#### 3.1. Search algorithm results

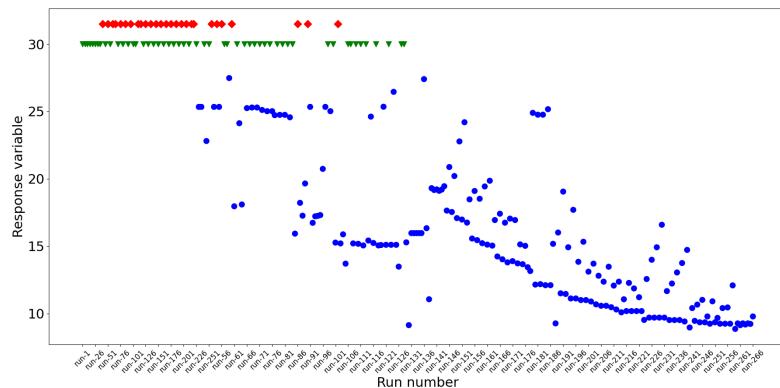
The search algorithm was allowed to search the space to determine the best material properties. The resulting response variables can be found in Figure 8. Where the blue circular markers indicate material datasets which developed a melt track and the red diamond markers did not have the energy density to produce a melt track. Due to the



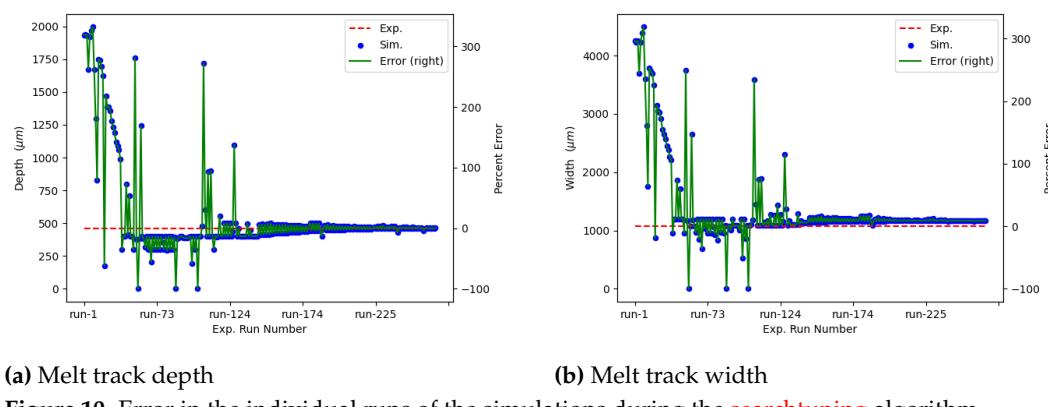
**Figure 8.** Response variable of the search algorithm for material properties and laser diameter where the blue circle marks developed a melt track and red diamonds did not.

vast difference in scales of the initial responses and the final response variables, a new plot was created which has a max Y value of just over 30. In this plot the blue circular markers are ones which have a response variable less than 30, the green triangle markers are those which completed with a melt track but had response variables greater than 30, and the red diamond markers are those which did not produce a melt track. In addition to these plots, the error in the width and depth were plotted and can be seen in Figure 10. In these plots, it can be seen that the error in the width is 8.83% and the error in the depth is 0.03%. It is not fully understood why all the error is coming from the width, however, it is theorized that this is a product of the difference in the size of the width vs the depth since the width is nearly triple that of the depth.<sup>6</sup>

<sup>6</sup> Updated figure captions to include marker explanations



**Figure 9.** Response variable of the search algorithm for material properties and laser diameter with max y axis value of 30 where the blue circle marks developed a melt track, green triangles developed a melt pool but had a response greater than 30, and red diamonds did not develop a melt pool.



**Figure 10.** Error in the individual runs of the simulations during the `searchtuning` algorithm

The search algorithm completed and reduced the combined error from over 600% when starting from the material properties found in the literature for the generic aluminum material properties, found in Table 6, to 9.1% when using the values found in Table 8.

**Table 8.** Optimized Aluminum material properties and laser diameter dataset for the developed simulation

Property	Material Temp.	Value
Laser absorption	880°C	16.8%
Laser absorption	922°C	10.0%
Thermal conductivity	922°C	32.2 $\frac{W}{mK}$
Thermal conductivity	1491°C	152.3 $\frac{W}{mK}$
Specific heat	733°C	2957.6 $\frac{J}{kgK}$
Laser diameter		0.864 mm

### 3.2. Search results validation

In order to ensure that the search algorithm results were valid across laser travel speeds and power levels, a range of 8 other parameters were compared with experimental results. These experiments and simulations were the setup and analyzed as was used for

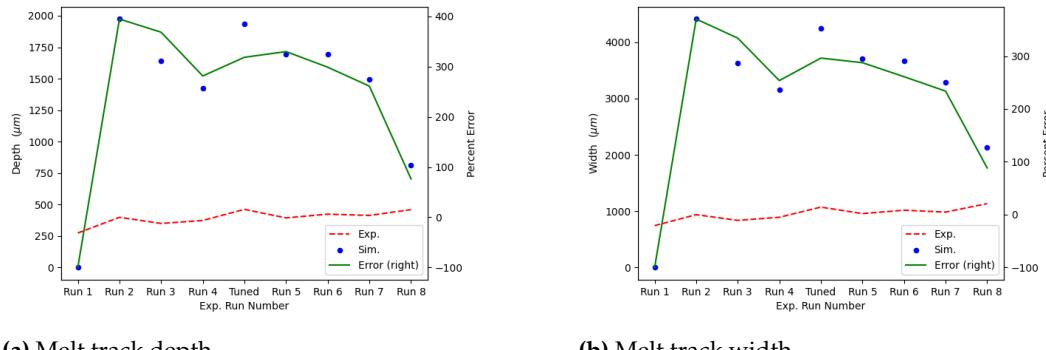
the [search algorithm tuning](#) experimental setup. The base experimental parameters can be seen in Table 7 with the scan speed and laser power being varied as seen in Table 9 and the simulation material properties were used from Table 8. The experimental result were

**Table 9.** Laser processing parameters validation [validation](#) processing parameters

Exp. Id.	Scan speed (mm/min)	Laser Power (W)
1	762	1000
2	762	1500
3	762	1250
4	1143	1250
5	1143	1500
6	1524	1750
7	1524	1500
8	1524	2000

collected and graphed in Figure 12 and 11 as the red dashed line and were used as the ground truth with which the models were compared.

These speeds and powers were first completed with the literature determined values from Table 6 and the results can be seen in Figure 11. Where the red dashed line is the

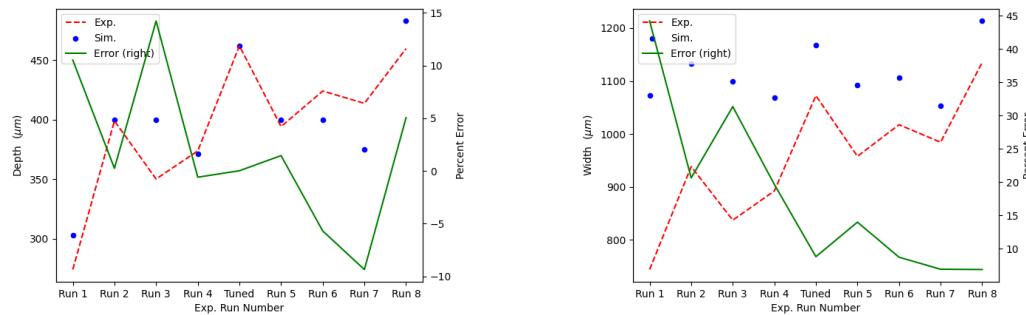


**Figure 11.** Comparison of experimental and simulated results for validation points with generic literature values for [aluminum](#) material dataset

experimental results, the blue dots are the simulation predictions, and the green line is the error when comparing the simulated results to the experimental results. These results show that over the 9 initial parameter sets, when a melt track was developed, the average absolute value of the error in the depth was approximately 290% and the average error in the width was approximately 265%. To put this into terms of the response variable of the search algorithm, the sum of the width and depth error, would be 555% combined error. Additionally, run 1 was unable to develop a melt pool, which is contrary to the experiments where all the parameter sets had a stable melt track. These results corroborate the results from Figure 8 which showed that the parameter set used for [the search algorithm](#)

tuning initial dataset of material properties found in the literature is wholly inadequate for simulating the process at hand.

In contrast to these results, the material dataset which was found in Table 8 was used to simulate each parameter set, and the results can be seen in Figure 12. Where the



(a) Melt track depth

(b) Melt track width

**Figure 12.** Comparison of experimental and simulated results for laser processing parameters validation points with optimized values for Aluminum material dataset

red dashed line is the experimental results, the blue dots are the simulation predictions, and the green line is the error when comparing the simulated results to the experimental results. These results show the average error in the width was approximately 17% and the average error in the depth was approximately 5%, which creates a combined error of only 22%. These results show that the optimized dataset is better at predicting the combined error of the simulation by over 500%. This results in a simulation which can be leveraged more intensely during the process development and build qualification process. These results are from a wide parameter set which encompasses most of the usable parameter space for deposition found experimentally. This improvement in accuracy will allow for a greater application in the model in the determination of an optimized parameter set for a given build. In addition, the model is seen to be more accurate at parameters above the parameter used during the optimization. This leads to the conclusion that if a more accurate simulation is needed, the optimization should be done at a parameter set near the desired parameter set.

#### 4. Discussion

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This model was developed in response to a need from those developing processing parameters and path plans for the metal AM process. They were needing a model which was capable of quickly simulating the heat flow and buildup for a range of path plans

<sup>7</sup> Added a discussion section to bridge the gap between the results and conclusions

and processing parameters. This will allow for researchers to more test a wider range of  
292 parameters and path plans in order to better find a global optimal as opposed to the local  
293 optimization which might be obtained with a smaller scale set of parameters or path plans,  
294 both experimental and simulated. With the current target of an efficient predictor of the  
295 thermal history completed, and an optimized input dataset for the high strength aluminum,  
296 it is possible to expand this model to predict more components of the build and increase its  
297 usefulness.  
298

The next target for the model is to add prediction of microstructure of the material.  
299 This is particularly important for the aluminum alloy investigated in this paper because  
300 it is the main driving factor as to the expected performance of the final part[28]. In order  
301 to predict the microstructural makeup of an AM build, it is necessary to first find the  
302 thermal history due to the tight linking between the thermal history, cooling rates, and  
303 microstructure which is developed. After this a range of models, both determination and  
304 probabilistic, can be used to predict the final microstructural makeup of a completed AM  
305 build[29]. These results can then be used to predict the final part performance. Moving  
306 forward, various methods of predicting the microstructure of the completed build will  
307 be investigated as an intermediate step to estimating the ultimate performance of the  
308 component.<sup>8</sup>  
309

## 5. ConclusionsDiscussion

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This work shows that the Nelder-Mead search algorithm is an appropriate multidimensional  
312 search algorithm for the determination of the optimal dataset for improved  
313 simulation accuracy. It was capable of reducing the simulated melt track depth and width  
314 of a set of processing parameters by over 500%, as shown in Figures 11 and Figure 12, which  
315 used datasets from literature (Table 6) and the optimized dataset (Table 8) respectively. This  
316 was done by defining the response variable (Equation 10) for the search algorithm to be  
317 the sum of the error in the width and the depth, which facilitated an efficient search. This  
318 methodology can be used to develop accurate simulations for any material which is not  
319 well published or to increase the accuracy of a simulation which utilizes approximations of  
320 first principles in order to increase its efficiency.  
321

<sup>8</sup> Added discussion about microstructure

<sup>9</sup> The discussion section was more properly renamed to the Conclusions

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