Mitigation of lack of fusion defects in powder bed fusion additive manufacturing

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A R T I C L E   I N F O

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A B S T R A C T

Components manufactured by additive manufacturing often exhibit improper fusion among layers and hatchies. Currently, there is no practical way to select process parameters and alloy systems based on scientific principles to mitigate these defects. Here, we develop, test and demonstrate a methodology to predict and prevent these defects based on a numerical heat transfer and fluid flow model for the laser powder bed fusion (PBF) additive manufacturing (AM). These defects are avoidable by adjusting laser power, scanning speed, layer thickness and hatch spacing. An easy to use parameter is proposed for practical use in shop floors. Relative susceptibilities of three widely used AM alloys are demonstrated using this parameter.

1. Introduction

In laser-assisted powder bed fusion (L-PBF) additive manufacturing (AM), components are made of multiple thin layers and hatchies. The soundness of the component depends on the fusion bonding among successive layers and hatchies [1,2]. Lack of fusion is detrimental to the mechanical properties of the component, and in extreme cases leads to part rejection [1,3]. Improper selection of laser power, scanning speed, laser spot radius, layer thickness, hatch spacing and alloy affect the formation of this defect [4]. Because of the involvement of many process parameters and alloys, currently there is no generally available methodology to guide engineers to avoid this defect.

Several attempts have been made to understand the effects of different process parameters on the lack of fusion defect for different alloy systems and significant volume of experimental data have been reported in the literature. Experimentally, lack of fusion defect was found to be reduced with increasing laser power for stainless steel [5], titanium alloys [4,6,7], aluminum alloys [8-10] and CoCrMo alloy [11]. Slower scanning speed was also found to reduce lack of fusion defect for titanium alloys [4,6,7,12,13] and aluminum alloys [8-10]. Thinner layer and small hatch spacing were proved to reduce this defect for stainless steel [5], titanium alloys [14] and aluminum alloys [8-10]. However, in these experiments, one process variable was varied while other parameters were kept constant. The brute force approach to evaluate a parameter space for avoiding such defects through empirical testing needs to be repeated for every alloy. Because of the involvement of many process parameters, experimental evaluation of the roles of process parameters on the prevalence of defects by trial and error is time consuming and expensive.

A recourse is to develop, test and validate a science based, easy to use methodology of avoiding these defects using heat transfer considerations. Heat conduction models [15,16] were used to calculate the molten pool shape and size based on which the extent of lack of fusion voids were estimated. However, these models neglect the effect of molten metal convection that is often the main mechanism of heat transfer inside the pool [1]. Neglecting molten metal convection results in inaccurate molten pool shape and size [17] and thus introduces uncertainty in the lack of fusion defects. In short, the existing powder bed fusion literature does not provide any rigorous methodology for selecting process parameters and alloy system to minimize lack of fusion defect. A quantitative understanding of the effects of process parameters and alloy properties on lack of fusion defect and a practical means to mitigate this problem based on scientific principles are needed but not currently available.

Here, we show, for the first time, how the lack of fusion defects during PBF can be minimized by using heat transfer and fluid flow calculations. To quantify the lack of fusion defect, we propose a dimensionless number that involves common process parameters such as laser power, scanning speed, spot radius, layer thickness and hatch spacing and different alloy properties such as density, thermal conductivity, specific heat and latent heat of fusion. The dimensionless number also includes molten pool dimensions calculated using a well-tested, three dimensional, transient heat transfer and fluid flow model of L-PBF process. This non-dimensional number provides a relative
scale to compare different alloys based on their susceptibility to the lack of fusion defect over a wide range of process parameters. The effects of all thermophysical properties of the alloys are considered during heat transfer and fluid flow calculations which is required for the lack of fusion number. Since the volume of the fusion zone depends on the processing parameters, thermophysical properties of the alloy and the geometry of the component, all of these factors have to be taken into account in the modeling to estimate the lack of fusion defect. In addition, the important non-dimensional numbers that are important in L-PBF such as Marangoni number and non-dimensional temperature are correlated with the occurrence of lack of fusion defect. Based on these findings we provide recommendations to mitigate lack of fusion defects in components. Although, the results presented here are for L-PBF process, the proposed lack of fusion number is applicable to all AM processes since it is formulated based on the underlying principles of the lack of fusion void formation that are similar in all AM processes.

2. Heat transfer and fluid flow model of PBF

A well-tested, three-dimensional, transient heat transfer and fluid flow model of L-PBF is used to calculate the temperature field and molten pool dimensions. The model solves the following equations of conservation of mass, momentum and energy [18–20]:

\[
\frac{\partial u_i}{\partial x_i} = 0 
\]

(1)

\[
\frac{\partial u_i}{\partial t} + \rho \frac{\partial (u_i u_j)}{\partial x_i} = \frac{\partial}{\partial x_i} \left( \mu \frac{\partial u_i}{\partial x_i} \right) + S_i 
\]

(2)

where \(\rho\) and \(\mu\) are the density and the viscosity of the alloy, respectively, \(u_i\) and \(u_j\) are the velocity components along the \(i\) and \(j\) directions, respectively, \(x_i\) is the distance along the \(i\) direction and \(S_i\) is the source term for \(i\)th momentum conservation equation. The energy conservation equation is written as [18–20]:

\[
\frac{\partial h}{\partial t} + \frac{\partial (u_i h)}{\partial x_i} = \frac{\partial}{\partial x_i} \left( \alpha \frac{\partial h}{\partial x_i} \right) - \frac{\partial (u_i \Delta H)}{\partial x_i} + S_h 
\]

(3)

where \(h\) denotes the sensible heat, \(t\) is the time, \(\alpha\) and \(\Delta H\) are the thermal diffusivity and the latent heat of fusion of the alloy, respectively, \(u_i\) is the velocity components along the \(i\) direction and \(S_h\) is the source term for the volumetric heat source and is represented as [1]:

\[
S_h = \frac{\xi \eta P}{\pi \tau t} \exp \left( -\frac{\xi (x_i^2 + y_j^2)}{r^2} \right) 
\]

(4)

where \(P\) is laser power, \(\tau\) is laser beam radius, \(\xi\) is power distribution factor varies between 1 and 3 [1], \(x_i\) and \(y_j\) are the distances from the laser beam axis along X and Y directions, respectively and \(t\) is the powder layer thickness.

The solution domain consists of substrate, deposited layers and hatches, powder bed, and shielding gas. The substrate material is same as the alloy powder. Convective and radiative boundary conditions are applied to all surfaces of the solution domain as [1]:

\[
- k \frac{dT}{dz} = \sigma (T^4 - T_s^4) + h_c (T - T_s) 
\]

(5)

where \(k\) is the thermal conductivity of alloy, \(\sigma\) is the Stefan–Boltzmann constant \((5.67 \times 10^{-8} \text{Wm}^{-2} \text{K}^{-4})\), \(\varepsilon\) is the emissivity, \(T_s\) is the ambient temperature and \(h_c\) is the convective heat transfer coefficient.

The convective flow of the molten metal is primarily driven by the surface tension variation on the top surface of the molten pool resulting from the spatial gradient of temperature. The resulting stress on the top surface of the molten pool, called Marangoni shear stress [20–22] along X and Y directions can be written as,

\[
\tau_x = \mu \frac{dx}{dz} = -\frac{dy}{dT} G_x 
\]

(6)

\[
\tau_y = \mu \frac{dy}{dz} = -\frac{dx}{dT} G_y 
\]

(7)

where \(\mu\) is the viscosity of the liquid metal, \(G_x\) and \(G_y\) are the two components of temperature gradient along X and Y directions and \(dy/dT\) represents the surface tension gradient with respect to temperature that has a negative value for most commonly used alloys that do not contain any surface active element [1]. Temperature dependent thermo-physical properties of the alloy powders [23] are provided in Table 1.

<table>
<thead>
<tr>
<th>Properties</th>
<th>SS 316</th>
<th>Ti-6Al-4V</th>
<th>AlSi10Mg</th>
</tr>
</thead>
<tbody>
<tr>
<td>Liquidus temperature (K)</td>
<td>1733</td>
<td>1928</td>
<td>867</td>
</tr>
<tr>
<td>Solidus temperature (K)</td>
<td>1693</td>
<td>1878</td>
<td>831</td>
</tr>
<tr>
<td>Thermal conductivity (W/m K)</td>
<td>11.82 + 1.06 \times 10^{-2} T</td>
<td>1.57 + 1.6 \times 10^{-2} T – 1 \times 10^{-6} T^2</td>
<td>113 + 1.06 \times 10^{-5} T</td>
</tr>
<tr>
<td>Specific heat (J/kg K)</td>
<td>330.9 + 0.563 T – 4.015 \times 10^{-4} T^2 + 9.465 \times 10^{-8} T^3</td>
<td>492.4 + 0.025 T – 4.18 \times 10^{-5} T^2</td>
<td>536.2 + 0.035 T</td>
</tr>
<tr>
<td>Density (kg/m^3)</td>
<td>7900</td>
<td>4000</td>
<td>2670</td>
</tr>
<tr>
<td>Latent heat of fusion (J/kg)</td>
<td>272 \times 10^3</td>
<td>284 \times 10^3</td>
<td>423 \times 10^3</td>
</tr>
<tr>
<td>Viscosity (kg/m s)</td>
<td>7 \times 10^{-3}</td>
<td>4 \times 10^{-3}</td>
<td>1.3 \times 10^{-3}</td>
</tr>
<tr>
<td>Absorption coefficient in liquid (\eta)</td>
<td>0.3</td>
<td>0.3</td>
<td>0.3</td>
</tr>
<tr>
<td>Absorption coefficient in powder (\eta)</td>
<td>0.7</td>
<td>0.7</td>
<td>0.7</td>
</tr>
<tr>
<td>Volumetric expansion coefficient (/K)</td>
<td>5.85 \times 10^{-5}</td>
<td>2.5 \times 10^{-5}</td>
<td>2.4 \times 10^{-5}</td>
</tr>
<tr>
<td>Young’s modulus (GPa)</td>
<td>206</td>
<td>110</td>
<td>68</td>
</tr>
</tbody>
</table>

Table 1: Thermo-physical properties of SS 316, Ti-6Al-4 V and AlSi10Mg [23]. Here ‘T’ represents temperature in K ranging from ambient to the solidus temperature.

More details about the implementation of this model for L-PBF are described in our previous publication [24] and are not repeated here. Only the salient features are indicated here. The conservation equations of mass, momentum and energy are discretized in the 3D Cartesian coordinate using the control volume method. All discretized equations are solved simultaneously to obtain enthalpy, velocity and pressure fields using the tri-diagonal matrix algorithm (TDMA) [25]. The temperature field is obtained from the enthalpy field by using temperature dependent specific heat of the alloy. The calculation procedure continues until all the hatches and layers are completed. These calculations are performed using an in-house Fortran code compiled using an Intel Fortran compiler. The process parameters used for the calculations are given in Table 2. The calculation time is approximately 5 h for a 20 mm long, 5 layers, 5 hatches build in a personal computer with a 3.40 GHz i7 processor and 8 GB RAM.

3. Results and discussions

Fig. 1(a) shows the three-dimensional temperature and velocity fields calculated using the heat transfer and fluid flow model during the building of first layer first hatch of a SS 316 build. The temperature and velocity fields on top (XY), transverse (YZ) and longitudinal (XZ) planes are shown in Fig. 1(b)–(d) respectively. The region bounded by the liquidus temperature isotherm (1733 K) of SS 316 represents the fusion
zone of the molten pool. The light blue region within the liquidus and the solidus temperature (1693 K) isotherms represents the mushy zone. Therefore, the 1693 K isotherm represents the molten pool boundary. The velocity vectors are represented by the black arrows whose magnitude can be found by comparing their length with the reference vector provided. The velocity vectors are radially outwards because molten metal flows from the high temperature to the low temperature. The laser beam travels in the direction of positive X-axis. Therefore, the molten pool is elongated in the opposite direction (negative X-axis). The molten pool exhibits a teardrop shape that is attributed to rapid scanning speed of the PBF process [1]. The track width is determined from the solidus isotherms on the transverse sections as shown in Fig. 1(d). Table 2

| Process parameters used for calculations. Packing efficiency is a measure of extent of voids in the powder bed and is adapted from the literature [24]. |
|---|---|
| Laser power, W | 60 |
| Scanning speed, mm/s | 250-1000 |
| Laser spot radius, mm | 0.050 |
| Layer thickness, mm | 0.030 |
| Hatch spacing, mm | 0.035 |
| Build length, mm | 20 |
| Substrate dimensions, mm × mm × mm | 22 × 5 × 2 |
| Packing efficiency | 0.5 |

Fig. 1. Temperature and velocity distributions for 1st hatch and 1st layer of a 20 mm long SS 316 build on a 24 mm long SS 316 substrate using 60 W laser power and 500 mm/s scanning speed on (a) 3D isometric section (b) top (c) longitudinal and (c) transverse planes. The length of the build is from x = 2.0 mm to 22.0 mm. Other process conditions are mentioned in Table 2. Scanning direction of the laser beam is along the positive x-axis.

Fig. 2. Comparison between the calculated and experimentally observed (a) width and (b) depth of the molten pool of a single layer single hatch builds of SS 316, Ti-6Al-4 V and AlSi10Mg at different linear heat inputs. The experimentally measured width and depth for SS 316 are adapted from Di et al. [5] and Li et al. [27], respectively. The experimental results for Ti-6Al-4 V and AlSi10Mg are taken from Gong et al. [12] and Kempen et al. [26], respectively. Gong et al. [12] and Kempen et al. [26] provided the macrograph from which Tang et al. [15] measured the dimensions.
evaluate lack of fusion defect.

Figs. 3(a–c) show the computed shapes and sizes of the molten pool transverse sections (YZ plane) for different hatches and layers of five hatches, three layers, builds of SS 316, Ti-6Al-4 V and AlSi10Mg, respectively. For the SS 316 and Ti-6Al-4 V builds, unmelted regions between the molten pools indicating improper fusional bonding among layers and hatches represent the lack of fusion voids. For the same processing conditions, molten pools in AlSi10Mg build are the largest due to its lowest density. Therefore, AlSi10Mg does not exhibit any lack of fusion voids at the processing condition considered. Figs. 4 (a–c) show the computed shapes and sizes of the molten pool transverse sections (YZ plane) for different hatches and layers of five hatches, three layers, builds of SS 316 at different heat inputs. Lack of fusion voids observed in Fig. 4 (a) are eliminated by increasing laser power and decreasing scanning speed, as shown in Fig. 4 (b) and (c), respectively.

From Fig. 4 it is clear that the lack of fusion defect depends on both laser power and scanning speed. Figs. 5 (a) and (b) show the relation of experimentally measured [5,11] lack of fusion void fraction with laser power and scanning speed, respectively. In these plots the experimental data are taken from the independent literature [5,11]. It has been found that amount of the lack of fusion voids is directly proportional to both the layer thickness and hatch spacing.

From Figures (3–6) it is clear that the lack of fusion defect depends on alloy properties and process parameters such as laser power, scanning speed, layer thickness and hatch spacing. However, other process parameters such as laser spot radius, absorptivity of the laser beam at the powder bed, molten pool width and depth and rate of heat transfer also govern lack of fusion defect in PBF [4,15]. Therefore, to quantify and provide better understanding of the effects of these governing factors on the lack of fusion defect in PBF, a non-dimensional lack of fusion number (L_F) is proposed and used here. This number, L_F consists of all important process parameters and alloy properties and is represented as:

\[ L_F = \frac{F (C_\text{w} + T + L) \int \left( \frac{h}{d} \right)^2 \, d}{\pi r^2} \]

All the symbols used in this equation are described along with their units and dimensions in Table 3. Eq. (8) clearly indicates that the lack of fusion defect is proportional directly to scanning speed, layer thickness and hatch spacing and inversely to laser power as shown in Fig. 5 and 6. Molten pool width and depth and Fourier number used in this equation are calculated using the heat transfer and fluid flow model. Recently,
we have shown how the molten pool dimensions vary with layers and hatches [24,29]. The dimensions do not change significantly after the 2nd hatch and the layer where the temperature field reaches steady state [24,29,30]. Therefore, molten pool dimensions used in Eq. (8) are taken at 2nd hatch after they reach the steady state. Since wider and deeper molten pool ensures proper fusional bonding among successive layers and hatches, both pool width and depth are in the denominator of Eq. (8). Other important material properties that affect pool dimensions and thus the lack of fusion defect such as viscosity and coefficient of surface tension are considered by including their effects during pool depth and width predictions. Fourier number indicates faster diffusive heat transfer relative to heat accumulation [31]. A high rate of heat transfer reduces the pool size and thus increases the lack of fusion defect. The term \( \rho C \Delta T L \) denotes the amount of heat needed to melt per unit volume of material. For a given heat input, an alloy with high energy required for melting exhibits smaller molten pool that increases the susceptibility of lack of fusion defect.

Fig. 7 shows that the experimentally measured [5,15,28,13] lack of fusion void fraction \( V_E \) for three commonly used alloys follows a linear relationship with corresponding \( L_F \), which is estimated using Eq. (8). A sample calculation for the estimation of the \( L_F \) for SS 316 is shown in the Supplementary document. Based on the trend of the data points presented in Fig. 7, the lack of fusion void fraction \( V_E \) can be expressed as:

\[
V_E = 15.3 L_F \tag{9}
\]

All statistical data to qualify this linear fit are summarized in Table 4. Eq. (9) is valid for the heat input range of 0.05–1.00 J/mm, which is widely used for major L-PBF applications [1]. In this equation, when \( L_F \) is zero there are no lack of fusion voids observed. To include a new alloy, the correlation in Eq. (9) needs to be updated by including new experimental data for that alloy in Fig. 7. However, the lack of fusion number is applicable for any alloy for a wide range of processing conditions.

The lack of fusion number, \( L_F \), provides a usable scale to estimate and compare the amount of lack of fusion void in L-PBF of different alloys. Fig. 8 compares three commonly used alloys based on their vulnerability to the lack of fusion defect at three different scanning speeds. SS 316 is the most susceptible to lack of fusion defect because of its smallest molten pool attributed to its relatively high density. Since rapid scanning reduces the pool size, amount of lack of fusion voids enhances with increasing scanning speed as shown in Fig. 8.

Lack of fusion defect depends largely on the molten pool shape and size governed by the flow of liquid metal driven primarily by the spatial gradient of interfacial tension, also known as the Marangoni stress [20–22]. Effects of Marangoni stress is quantified by the Marangoni number [31,32]:

\[
Ma = \frac{\gamma}{\mu \alpha} \frac{d \delta}{dT} \frac{\Delta T}{\Delta Y} \tag{10}
\]

where \( \mu \) is the dynamic viscosity, \( \alpha \) is the thermal diffusivity of the
alloy, $\delta$ is the characteristic length of the molten pool, which is taken as the width of the molten pool, $\Delta T$ is the difference between the peak temperature inside the pool and the solidus temperature of an alloy, and $\frac{d\gamma}{dT}$ is the surface tension gradient with respect to temperature. For most alloys that do not contain any surface active elements, this quantity is negative [1]. The peak temperature and pool width required for the calculations are estimated using the heat transfer and fluid flow model. Higher Marangoni number indicates vigorous flow of liquid metal inside the pool that increases the molten pool width and ensures proper fusional bonding among successive layers and hatches. Therefore, lack of fusion defect decreases when hatch spacing and layer thickness are constant for processes with higher Marangoni number as shown in Fig. 9.

Assuming a constant cross section, hatch spacing, and layer thickness, higher molten pool peak temperature may indicate heat accumulation and consequently a larger molten pool which facilitates better bonding of the depositing metal with the previously deposited metal. Therefore, monitoring of peak temperature during AM can be used as an indicator of the extent of lack of fusion defect. A non-dimensional temperature $T^*$ can reveal the change in the amount of lack of fusion voids due to rise in peak temperature [32]:

$$T^* = \frac{T_P}{T_L}$$  

where $T_P$ and $T_L$ are the peak temperature and liquidus temperature of
the alloy, respectively. The peak temperature required for the calculations may be either estimated from the heat transfer and fluid flow model or determined with a thermo-camera. Fig. 10 shows that the lack of fusion defect decreases with an increase in peak temperature. The peak temperatures for the parameter range considered here are below the boiling point of the alloys and keyholes do not form during the process. However, instabilities in the keyholes formed at very high power density may result in porosity that are different from the lack of fusion voids described here [1,33]. Fig. 9 and 10 show that the lack of fusion voids can be effectively minimized by enhancing Marangoni number and non-dimensional temperature during the process by adjusting different processing conditions.

4. Summary and conclusions

A well-tested, three-dimensional, transient heat transfer and fluid flow model is used to calculate the temperature and pool dimensions used for the dimensionless calculations. The results presented here provide, for the first time, a quantitative basis for minimizing lack of fusion defect in components made by L-PBF based on scientific principles. Below are the specific findings.

1) A non-dimensional number that considers AM process parameters and alloy properties is developed and tested for preventing lack of fusion defects for a wide range of process conditions and three commonly used alloys.
2) Among the three commonly used AM alloys considered, the biggest molten pool of AlSi10Mg for a given set of AM process parameters minimizes lack of fusion defect. Under the same processing conditions, AlSi10Mg and SS 316 are the least and most vulnerable to the
lack of fusion defect.  
3) A larger heat input per unit length obtained by reduction of scanning speed or an increase in laser power or both results is larger liquid pool and lower occurrence of lack of fusion defect.  
4) A high value of Marangoni number that indicates vigorous circulation of the liquid metal inside the molten pool correlated well with the reduction of the lack of fusion defects.  
5) High values of peak temperature also correlated well with the reduction of the occurrence of lack of fusion defects. Since the temperature can be monitored during deposition, this correlation can be used to reduce lack of fusion defects.  

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Appendix A. Supplementary data  
Supplementary material related to this article can be found, in the online version, at https://doi.org/10.1016/j.jmapro.2018.10.028.  

References  


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