

Laser powder bed fusion of AlSi10Mg alloy: Numerical investigation on the temperature field evolution

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ABSTRACT

Additive manufactured parts are subjected to intense thermal gradients and high temperature peaks which affect mechanical properties. Such thermal cycles can cause distortions, residual stresses and microstructural heterogeneities. Since the experimental measurement of the temperature field is extremely difficult, numerical simulation can be used to obtain a description of the phenomenon. Here, a three-dimensional computational model for the prediction of the temperature field during the laser powder bed fusion process on AlSi10Mg alloy was developed. Scan path, the geometry of the heat source and the progressive generation of the part during the process have been simulated with finite element method. This approach was used in a small scale representation, as the extremely fast temperature gradients, high scanning speeds and amount of thermal energy input make the phenomenon extremely localized. The predicted melt pool size, compared with microstructural analysis results on reference samples, was used to validate the computational model.

1. Introduction

Nowadays, the production of near-net-shape metal components via Additive Manufacturing (AM) techniques is gaining increasing attention from both industrial sectors and researchers due to high design versatility [1]. However, the quality of additively manufactured parts is not always up to the industrial standards defined for traditional manufacturing methods. One of the main challenges for AM acceptability is the definition of rigorous certification methodologies balanced with the unique demands and features of AM [2]. The definition of predictive tools and standardized testing protocols is fundamental to ensure process reliability and reproducibility [3]. From this point of view, a detailed description of the several phenomena occurring during the manufacturing process is the key to understanding the linkage between the process and the microstructural features or the overall mechanical response of printed parts. The complex heat input dynamics causes local thermal histories characterized by severe gradients and cooling rates which are responsible for residual stresses and strains, defect formation mechanisms and of different microstructural morphologies. Thus, the part quality is strictly connected to the temperature field, which is the result of the process features and its parameters [2,4] *In situ* measurements of the thermal profile are quite challenging due to the high spatial and temporal resolution required [4]. Contact sensors

(thermocouples) [5,6] were used to monitor the temperature change of substrate points during the deposition. Although they have high accuracy and they are easy to calibrate, their major weakness is the inability to describe the thermal field along the energy input source path. Contactless measurements with pyrometers or thermal image cameras are more suitable for AM process monitoring since they can provide information about the area with the highest thermal dynamics. Nevertheless, multiple phenomena overlap during deposition, such as partial absorption and reflection of laser power, phase changes, evaporation, material expulsion and Marangoni flow may cause a local variation of the optical properties of the surface that is quite difficult to evaluate [7]. These technologies require an accurate estimation of the surface emissivity and the chamber transmissivity [8]. Researchers have proposed several techniques to correctly evaluate these parameters making the experimental set-up particularly complex [9], as highlighted by Yan et al. [4]. To overcome these limits, there has been a continuous effort in the development of numerical models aimed to simulate AM processes and to obtain information about the thermal field. Li et al. investigated the relationship between process parameters, the thermal profile (peak temperature, molten pool lifetime) and defects formation during a selective laser melting process of commercially pure Ti powders, comparing the numerical results with experimental microstructural analysis [10]. Wei et al. [11] proposed a 3D FEA model to evaluate the

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Table 1
Process Parameters.

Parameter	UoM	Value
Layer thickness	mm	0.05
Laser power	W	350
Scanning speed	mm/s	1150
Hatch distance	mm	0.17
Base temperature	°C	150
Initial powder temperature	°C	65
Recoat time	s	~2.5

influence of process parameters, such as laser power and scanning speed, on the thermodynamic behavior of the melt pool. Their results highlighted a strong Marangoni flow dependence on process parameters and that an unstable flow is responsible for defects generations within the printed material. Li et al. [12] developed a thermo-mechanical model to investigate the main features of a “point-exposure” scanning strategy and, from a multi-track simulation, they observed that the highest residual stresses appeared at the overlap of different tracks. Fergani et al. [13] implemented an analytical model in Matlab to predict both the temperature field and the residual stresses and strains maps during a selective laser melting process on 316 L alloy. They found that the maximum tensile residual stresses were along the building direction and concluded that selecting an adequate scanning path may result in an optimal residual stress profile for the component. The information on the transient temperature field has also been used to predict the grain structure, coupling the thermal analysis with Cellular automata [14] or Phase-Field method [15], or to predict the phase distribution, implementing phase evolution criteria [16,17]. Therefore, the prediction of the thermal field evolution during the deposition process has a very strong capability as it is useful to understand AM process dynamics, to predict final part key features or to optimize process parameters without extensive experimental investigations. In this study, a thermal transient computational model for the description of the thermal evolution during a laser powder bed fusion LPBF process on AlSi10Mg alloy is presented and validated. The model was developed with the commercial FEM code MSC MARC v2020. The main features of the thermal field were compared to the results of a metallographic analysis on reference samples.

2. Materials and methods

2.1. Sample fabrication and material properties

AlSi10Mg samples were produced using a SLM280 Twin machine, with two parallel IPG fiber lasers and a $280 \times 280 \times 365 \text{ mm}^3$ build envelope, in a protective Argon atmosphere. The maximum laser scanning speed is 10 m/s. The laser beam follows a raster pattern with a 67° change of orientation between consecutive layers in order to guarantee a high densification level of the final part. The metallographic analysis was carried out on vertical as-built cylindrical specimens whose dimensions are 8 mm (diameter) x 36 mm (height). Process parameters are listed in Table 1. The powder initial apparent density, measured according to ASTM-B417, is 1.52 g/cm^3 .

The accurate estimation of the thermo-physical temperature-dependent properties of the selected alloy is mandatory for a thorough material modeling. Indeed, during AM process, the material undergoes a complex thermal history over a wide temperature range, multiple melting and solidification. Criales et al. [18] underlined the strong dependence of the temperature profile on material properties such as conductivity, density and specific heat. In this study, those properties were obtained from *Sente JMatPro* simulation software given the initial powder chemical composition, estimated according to ASTM E1479. The effect of the latent heat was modeled through a significant discontinuity of the specific heat. The powder bed conductivity might be two times lower than the bulk material. It depends on the powder porosity, on the size

and distribution of the particles, and the gas filling the voids. The powder conductivity was estimated with a semi-empirical model [19]:

$$k_p = \frac{k_s(1 - \varepsilon)}{1 + \psi \left(\frac{k_s}{k_p} \right)} \quad (1)$$

where k_p , k_s and k_g are powder, bulk material, and gas conductivity, ψ is an empirical coefficient and ε is the porosity:

$$\varepsilon = \frac{\rho_s - \rho_p}{\rho_s} \quad (2)$$

ρ_s and ρ_p are the powder and solid density.

2.2. Heat transfer in LPBF

The distribution of the temperature T in a 3D cartesian domain satisfies the following equation,

$$\rho c_p \frac{\partial T}{\partial t} = \frac{\partial}{\partial x} \left(k(T) \frac{\partial T}{\partial x} \right) + \frac{\partial}{\partial y} \left(k(T) \frac{\partial T}{\partial y} \right) + \frac{\partial}{\partial z} \left(k(T) \frac{\partial T}{\partial z} \right) + Q \quad (3)$$

In which ρ , c_p and k are the material density, specific heat and conductivity and Q is the volumetric heat input. The initial temperature distribution is expressed as,

$$T(x, y, z) = T_0, t = 0 \quad (4)$$

Where T_0 is the initial powder temperature of 65 °C. During the deposition, the laser beam travels upon the powder bed, melting the underlying layer and allowing the formation of a metallurgical bond with the already deposited tracks. At the same time, heat transfer mechanisms such as conduction, convection and radiation play a crucial role on the surface, according to:

$$k \frac{\partial T}{\partial n} = q - q_{conv} - q_{rad} \quad (5)$$

q is the heat input flux. The heat dissipated by conduction is much higher than the one dissipated by convection, q_{conv} , and radiation, q_{rad} [20]. Due to the small scale of this phenomenon and the very high temperature gradients, conduction with the powder bed and the underlying layer is dominant [21,22]. To decrease the computational complexity of the analysis, the heat loss through the powder bed was modeled by an equivalent convective flux, according Li et al. [23]. The heat loss to the Argon environment was neglected. Radiative heat transfer is more effective than convection within the melt pool but still less than conduction [21,22]. Thus, a constant value of 0.4 [7,17] was chosen for the material's emissivity. The amount of laser energy absorbed by the material is difficult to evaluate given its high dependence on material thermo-physical properties, surface roughness, beam wavelength, and laser distance and orientation from the surface. The choice of a constant absorptivity coefficient has been widely used in literature [10,11,25,26]. In this work, an absorptivity coefficient of 0.3 was selected after an inverse calibration process aimed to reduce the error between the experimental and the predicted melt pool size, as estimated in [24].

2.3. Computational model

The multiphysics phenomena that occur during the process have led researchers to carry out modeling assumptions and simplifications in order to obtain a trade off between an accurate solution and the model complexity. Surely, one of the main problem in AM process simulation is the extensive computational time required to represent accurately the heat input phenomenon and its extremely rapid kinetics. As highlighted by Liu et al. [27], the temporal and spatial discretization should be detailed enough to ensure a continuous laser movement in the numerical simulation. To obtain information about the thermal field evolution at a

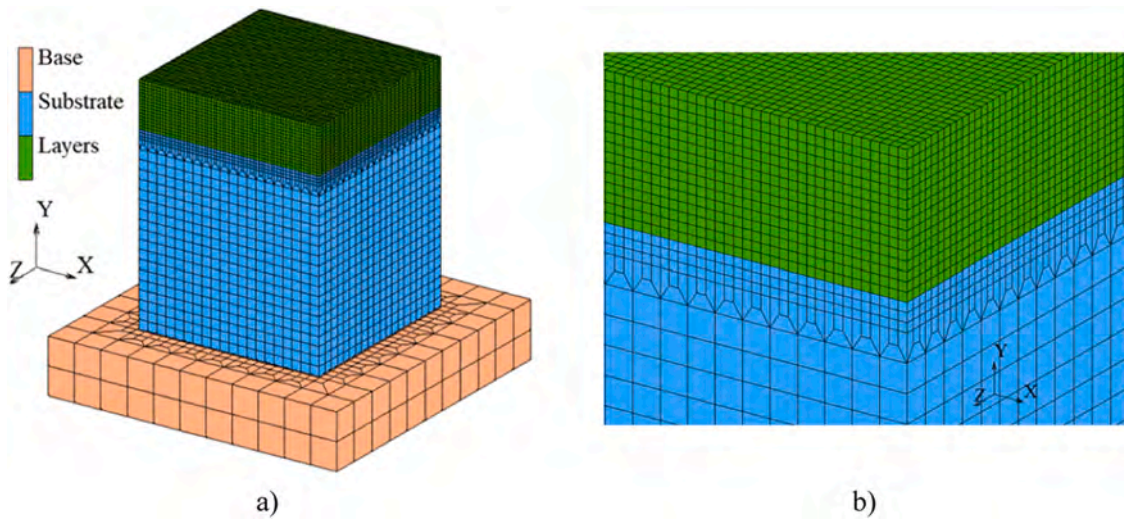


Fig. 1. Numerical model and computational grid.

part-scale level, the scale gap between the part dimension and laser spot size and the discrepancy between the time step required and the whole part generation time would made the analysis time prohibitive. This awareness led to the development of accurate simulation strategies to reduce the computational cost without losing accuracy for part-scale models. For instance, Ghanbari et al. [28] have subdivided the calculus domain in two scales, local and global. A small adaptive-local model with a very fine spatial and temporal discretization solved the thermal

field around the laser heat input, while a global simulation with larger grid size and time step calculated the temperature evolution far from the area with the highest kinetics. The two temperature profiles were then combined together. Liu et al. [27] developed and implemented in a FEM code a characteristic time-based heat input (CTI) model in which the energy input phase was simulated in two numerical steps. The scanning time of a track is divided in a heating time, in which an integrated energy input is applied, and a cooling phase without energy input. Other

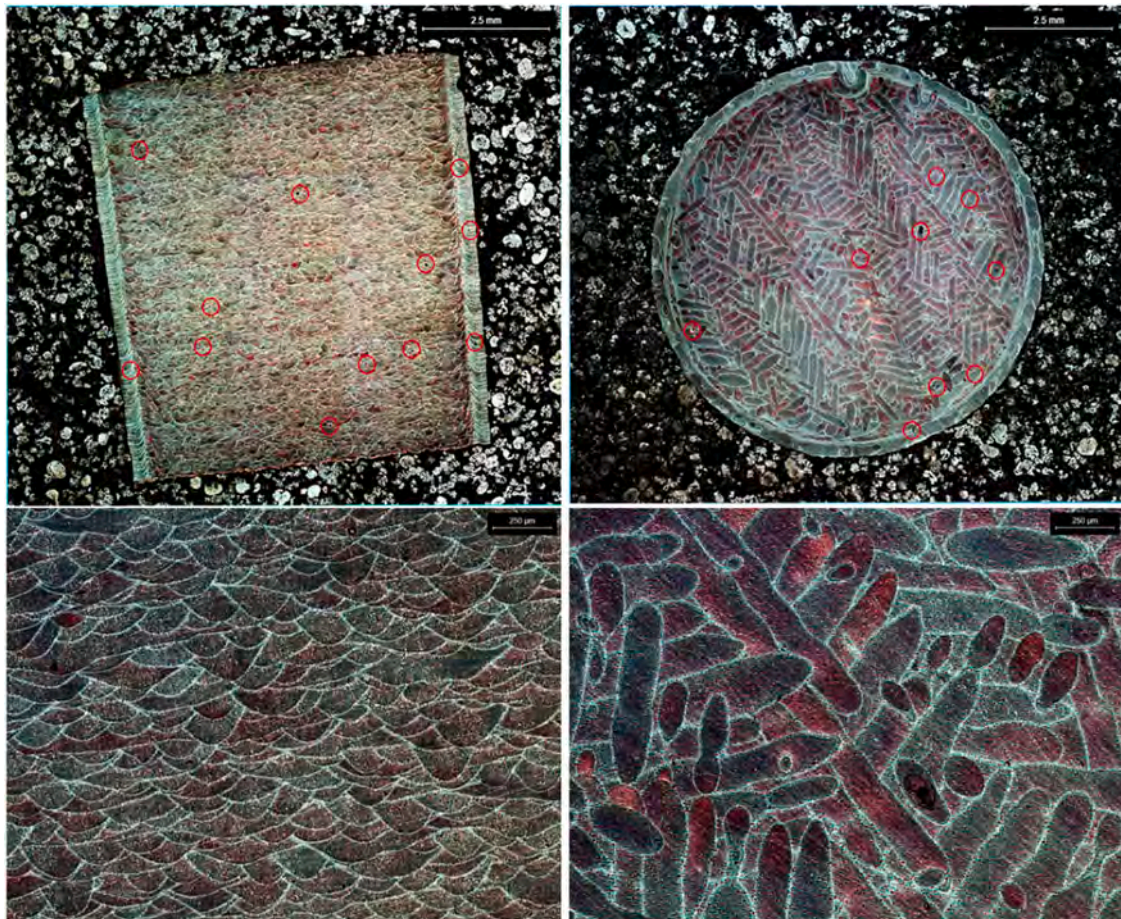


Fig. 2. Optical microscopy images - cross-section (left) top view (right) at different magnification levels.

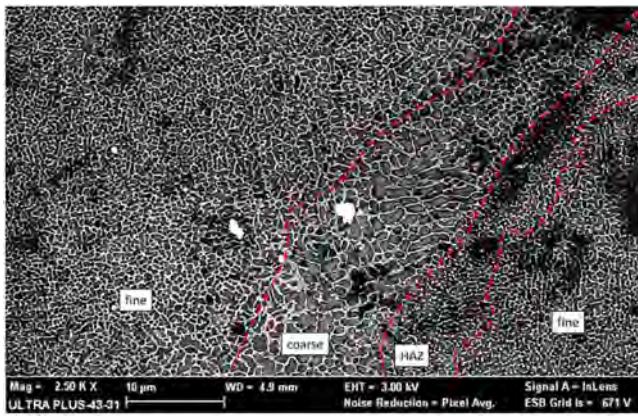


Fig. 3. SEM image shows the typical cellular microstructure of the melt pool cross-section.

authors [29–31] proposed a superposition-based semi-analytical model in which an analytical solution for point sources in a semi-infinite space is combined with numerical fields to impose boundary conditions. At the same time, the thorough description of the melt pool behavior and of the various phenomena that occur at the melt pool level requires quite accurate models that closely follow the movement of the laser input source [32]. Indeed, meso-scale models have been widely used in literature to describe local phenomena at the melt pool level to investigate key microstructural features [14,26,33] or the influence of process parameters on the thermal behavior or on defect formation mechanism [10,11,25]. This study is aimed to correlate the microstructural evidence to the thermal evolution the material undergoes during deposition, thus a meso-scale approach was chosen. Fig. 1 illustrates the computational grid, composed of hexahedral elements, with 8 integration points. The powder bed base is $1.6 \times 1.6 \times 0.3 \text{ mm}^3$, with a grid size of $0.15 \times 0.15 \times 0.15 \text{ mm}^3$. The specimen is modeled with a $1 \times 1 \times 1.25 \text{ mm}^3$ block, with a finer mesh of $0.016 \times 0.016 \times 0.016 \text{ mm}^3$ in the deposited layers. A coarser grid of $0.05 \times 0.05 \times 0.05 \text{ mm}^3$ was adopted far from the laser scan area. A glue contact was selected to replicate the welding link formed in LPBF process between the base and the part. The time step was chosen to guarantee a continuous laser motion upon the powder bed during deposition, so a constant value of $1.5 \times 10^{-5} \text{ s}$ was selected. During the recoat time, the time step gradually increases to a maximum of 0.15 s. The experimental parameters reported in Table 1 were used in the simulation.

The powder bed was modeled as a homogeneous mean [10,25] with scaled-down thermo-physical properties and only a thermal analysis was conducted [10,12,33], neglecting the melt pool fluidity due to its limited size [25]. To describe the change of conductivity between powder and solid, a hybrid quite-inactive element method was used. At the beginning of the analysis, all layers were inactive and not included in the calculus. Upon activation, the elements were in a quiet state, with a scaled-down conductivity, until they were crossed by the laser beam. The complex heat addition phenomenon was simulated by a spatially variable thermal flux along a user-defined path. In literature, various solutions have been employed to model the laser beam, such as the Gaussian heat source [11,28], Goldak's double ellipsoid [27,34], or a cylindrical geometry [35], while some authors proposed their own heat source models [36,37]. In this study, the laser input source was modeled using the weld flux tool with a cylindrical-shaped geometry, which allows a more focused heat input and more penetration in the material than Goldak's ellipsoid. The radius of the cylinder is the same of the laser beam input.

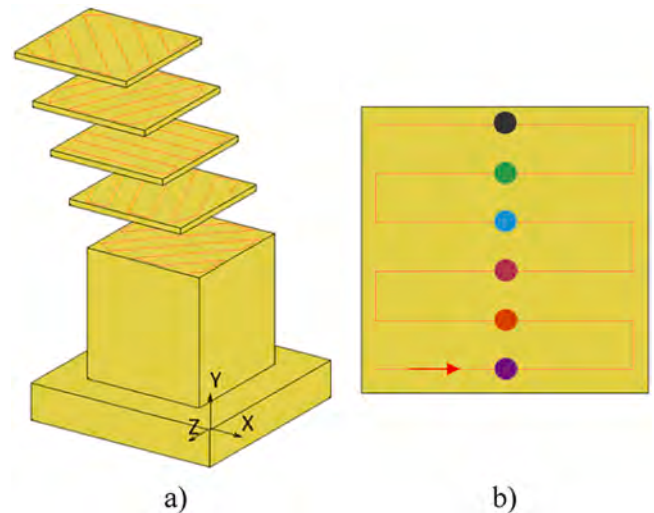


Fig. 4. (a) Scanning path and (b) measure points for the third layer.

3. Results and discussion

3.1. Sample characterization

As-built LPBF AlSi10Mg macro- and microstructural features are strongly dependent on the manufacturing process [38]. Given the high anisotropy of AM parts, vertical section and cross-section surfaces were extracted and analyzed. All resin-embedded samples were polished and etched by Keller's reagent. The microstructural features were observed using LEICA DM 6000 M optical microscopy (OM) and ZEISS ultra plus scanning electron microscopy (SEM). The optical microscopy images, presented in Fig. 2, highlight a strong texturization of the metallic matrix, related to the layer-by-layer nature of LPBF. Indeed, the vertical plane shows the typical semicircular morphology of the overlapped melt pools and in the cross-section plane, the laser scan tracks are identifiable. To estimate the melt pool size, these images were analyzed by ImageJ software; the melt pool depth was determined to be $110 \pm 20 \mu\text{m}$ and the width $220 \pm 30 \mu\text{m}$. The macrostructure along the cross-section and vertical planes is not uniform due to a 67° rotation of the scanning path for consecutive layers. As shown in Fig. 3, the major microstructural heterogeneities were found within the melt. The core of the melt pool exhibits a fine cellular structure of the $\alpha\text{-Al}$ phase surrounded by a continuous network of eutectic Al-Si precipitates. The cell size becomes coarser near the melt pool borders due to a change of solidification parameters [39]. Similarly to welding processes, outside the melt pool profile, there is a Heat Affected Zone (HAZ) in which Si phase coarsening causes a local rupture of the eutectic network.

3.2. Single-Layer analysis

A multi-track, multi-layer simulation was performed and the results were compared to the microstructural evidences. Fig. 4a illustrates the five reference layers with their corresponding scanning path. Firstly, the temperature evolution developed during the deposition of the 3rd layer (Fig. 4b) was investigated.

Fig. 5 illustrates the transient temperature distribution at different positions of the laser beam input. During the deposition of the first track (Fig. 5a), the influence of the heat input is localized mainly in the scanned area, in which the temperature increase is significant. On the contrary, the temperature of the remaining part of the powder bed is still close to the initial temperature of 65°C . Due to the discontinuity of thermal conductivity between the powder and the bulk material, the heat transmission by conduction towards the powder is slower than the laser movement. In fact, the front isotherms, that face the unscanned

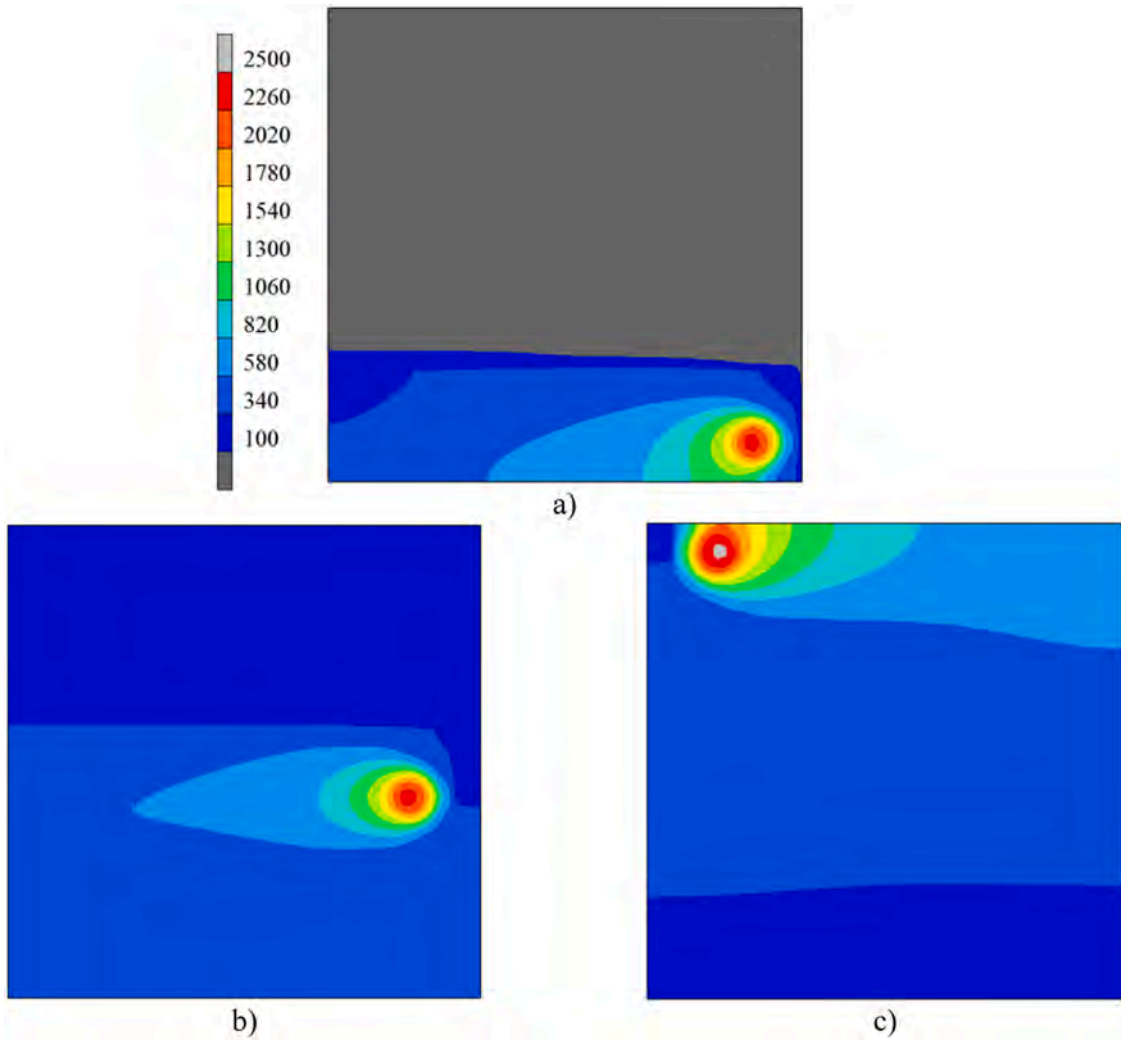


Fig. 5. Top view of the temperature distribution during the deposition of the third layer (a) end of 1st track (time 0.73 ms) (b) end of 3rd track (time 2.52 ms) (c) end of 6th track (time 5.31 ms).

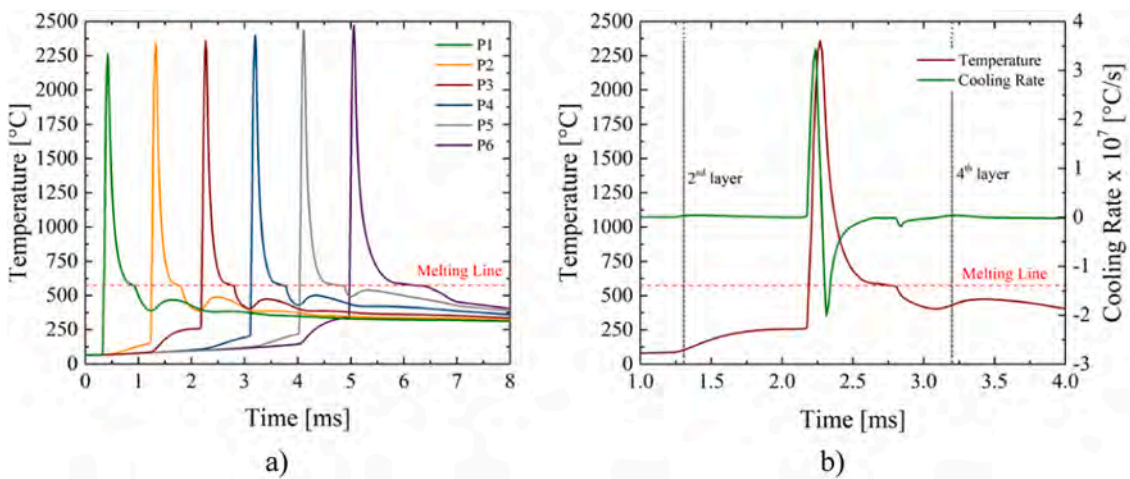


Fig. 6. a) Temperature evolution at the center of consecutive scanning tracks, b) temperature profile and the cooling rate at the center of the third scanning track.

zone, are denser than those at the back of the laser path where the material is in a solid-state state. Even during the deposition of the third track (Fig. 5b), the thermal distribution around the laser beam exhibits an asymmetrical behavior, but the powder bed temperature starts to

increase because of the longer interaction time with the laser beam. This process of heat accumulation starts to decrease during the deposition of the last track (Fig. 5c), where the material starts to cool down.

The thermal evolution of the center of each track, presented in

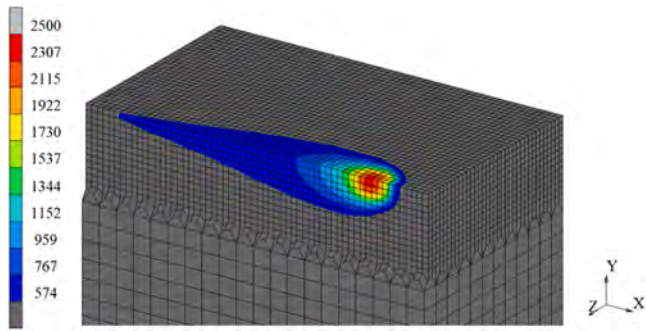


Fig. 7. Cross section view of the numerical temperature profile of the melt pool.

Fig. 6a, is characterized by very high cooling rates and temperature peaks, which far exceed AlSi10Mg melting temperature (574 °C). These results are in good agreement with the results obtained by Du et al. [25] for similar process parameters on the same alloy. The temperature peaks increase gradually from the first track to the last one, due to the heat accumulation during the deposition in the material. The influence between subsequent tracks can be clearly seen. For instance, the temperature at the center of the third track starts to gradually increase during the deposition of the second track. A content temperature increase during the deposition of the fourth track is also appreciable. A detail of the temperature change and cooling rate for the center point of the 3rd track is illustrated in Fig. 6b. An accurate description of the cooling profile and the phase change between liquidus and solidus temperatures is obtained taking into account both latent heat and material properties' temperature-dependence. For this temperature profile, the numerical cooling rate between the liquidus and solidus temperatures is $2.45 \times 10^5 \text{ s}^{-1}$, which is accountable for the very fine cellular structure observed in SEM images.

3.3. Melt pool analysis and model validation

In Fig. 7, a detail of the numerical melt pool profile during the laser motion is presented. The predicted melt pool dimensions were measured at the maximum extension of the thermal profile and they were estimated, as shown in Fig. 8, with the intersection of the temperature profile and the melting line. The results are in good agreement with the experimental data presented with their corresponding uncertainty in Fig. 9. The error in the prediction of both the melt pool width and depth is less than 2%.

The thermal histories within the melt pool were extracted in the five

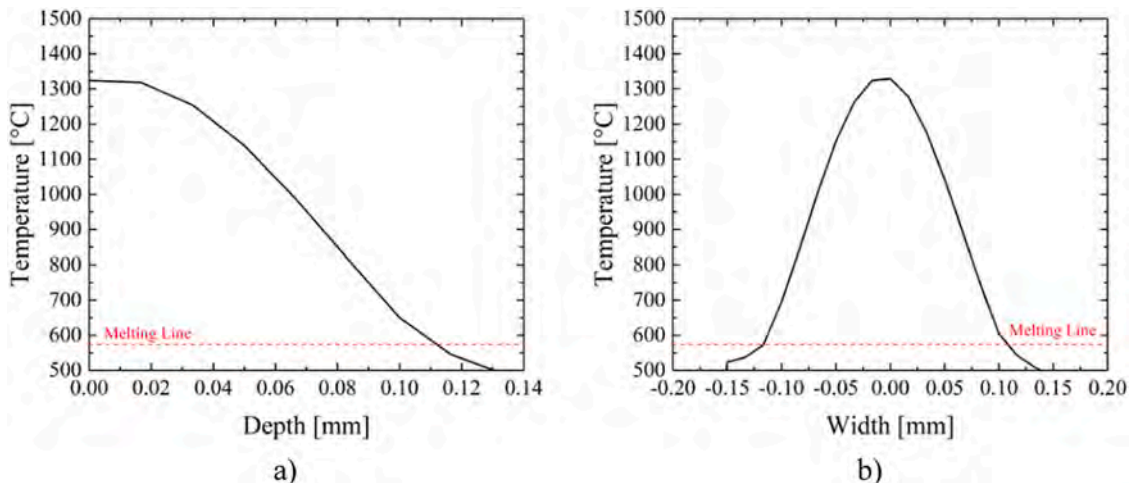


Fig. 8. Thermal profile extracted (a) along the depth (b) along the width of the melt pool.

measure points illustrated in Fig. 10.

The temperature profiles change meaningfully between the melt pool core and periphery. The latter has a more controlled thermal evolution, with lower peaks and slower cooling rates, Fig. 11, either along the width and the depth. This is in agreement with the grain size heterogeneity presented in the metallographic analysis, Fig. 3. The cell dimension is strongly dependent on the cooling rates. Along the width, the point at the periphery (W3) undergoes re-melting during the deposition of the second and the third track, as it can be clearly seen in

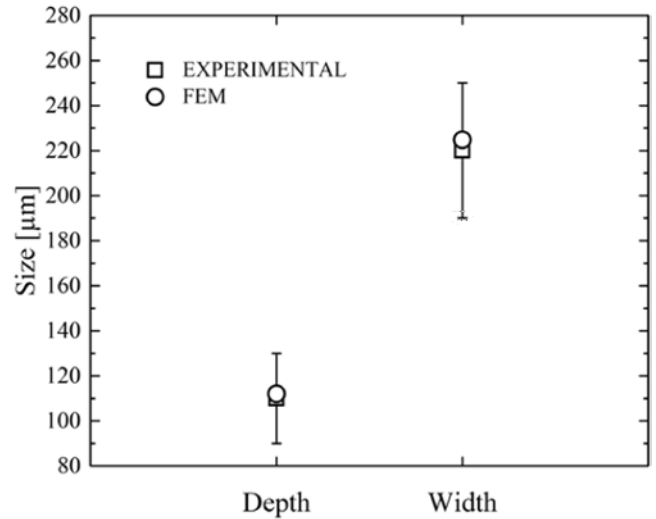


Fig. 9. Comparison of experimental and predicted melt pool size.

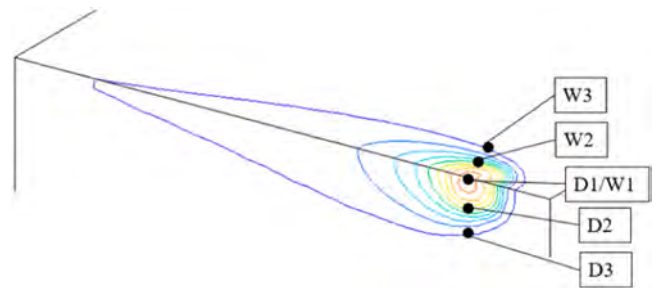


Fig. 10. Measure points within the melt pool.

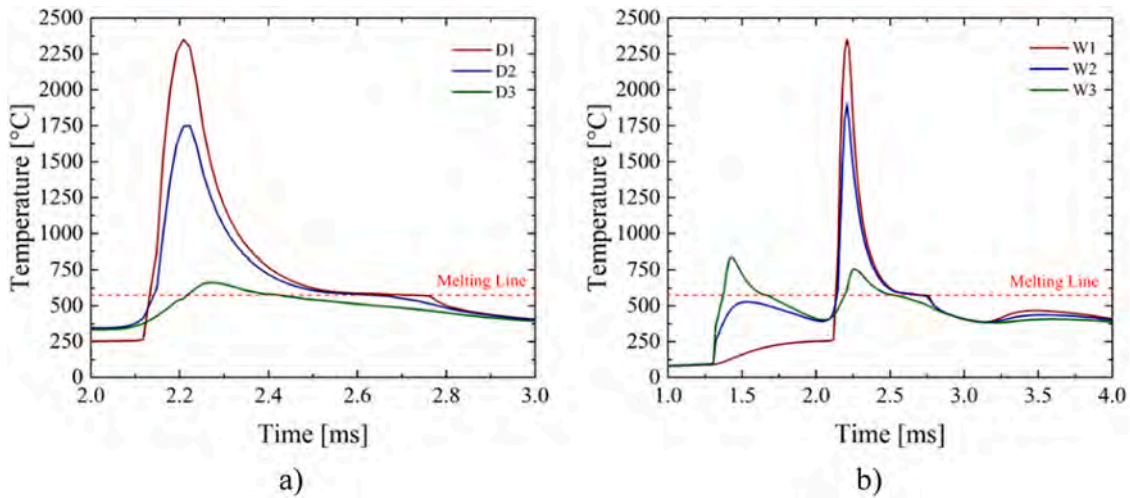


Fig. 11. Temperature distribution profiles within the melt pool (a) depth and (b) width.

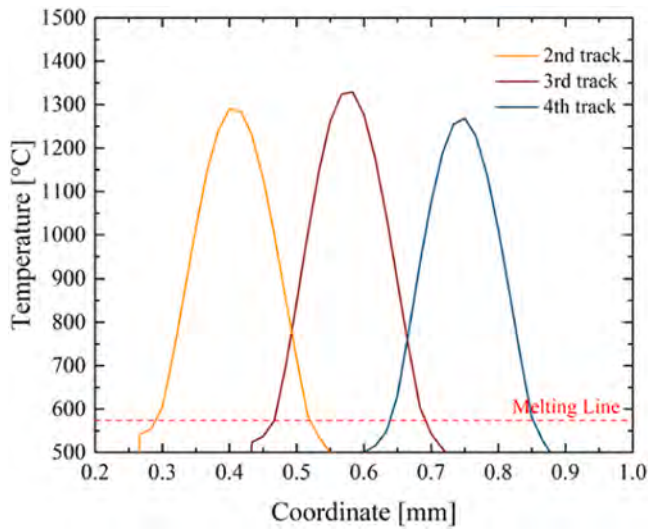


Fig. 12. Overlap of temperature profiles for subsequent tracks.

Fig. 11b. Indeed, the measured melt pool width is greater than the hatch spacing (0.17 mm), and this condition must be satisfied in order to guarantee a high densification level and reduce voids formation after the

scans. The overlap of the temperature profiles for three consecutive tracks is shown in Fig. 12. The melt pool width increases slightly during the deposition for the heat accumulation phenomenon. The predicted melt pool width is 213.4, 224.8 and 229.5 μm during the deposition of the second, third and fourth track respectively.

3.4. Multi-layer analysis

During a LPBF process, the material undergoes a complex thermal cycle, with repeated melting and solidification. The deposition of a new layer requires an additional energy input that has a variable thermal influence on the solidified layers. Figs. 13 and 14 show the thermal cycle measured at the center point of the first and second layers. Each layer undergoes severe re-melting during the deposition of the subsequent one: this condition guarantees the formation of a metallurgical bond between the two. The experimental melt pool size is a little greater than the sum of two layers thickness. Since the phenomenon is extremely localized, the thermal influence on other layers is limited, as supported by the numerical results. Fig. 15 underlines a self-similarity between the thermal histories of corresponding points of different layers. This result is corroborated by the experimental evidence since no microstructural gradients were found along the building direction.

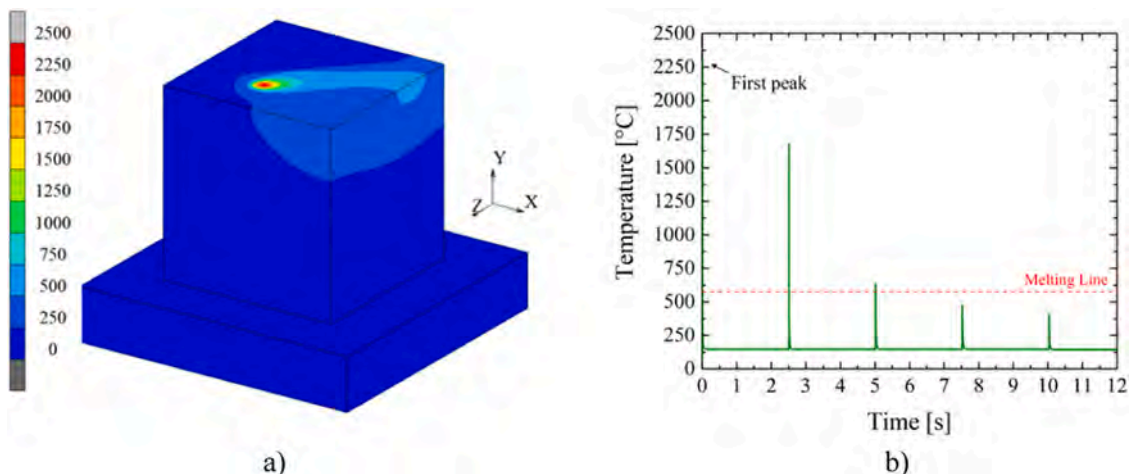


Fig. 13. Temperature map and thermal cycle evolution for the point at the middle of the 1st layer.

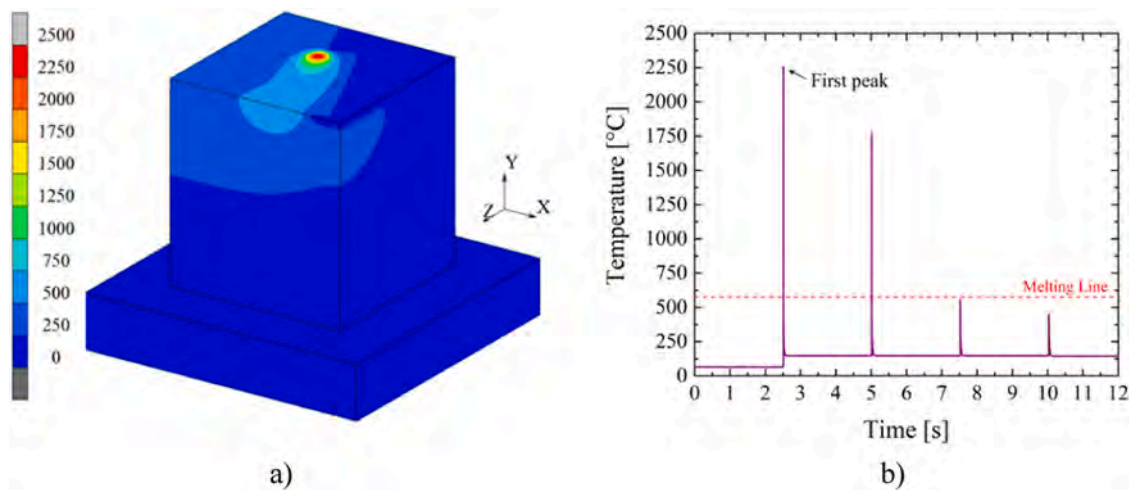


Fig. 14. Temperature map and thermal cycle evolution for the point at the middle of the 2nd layer.

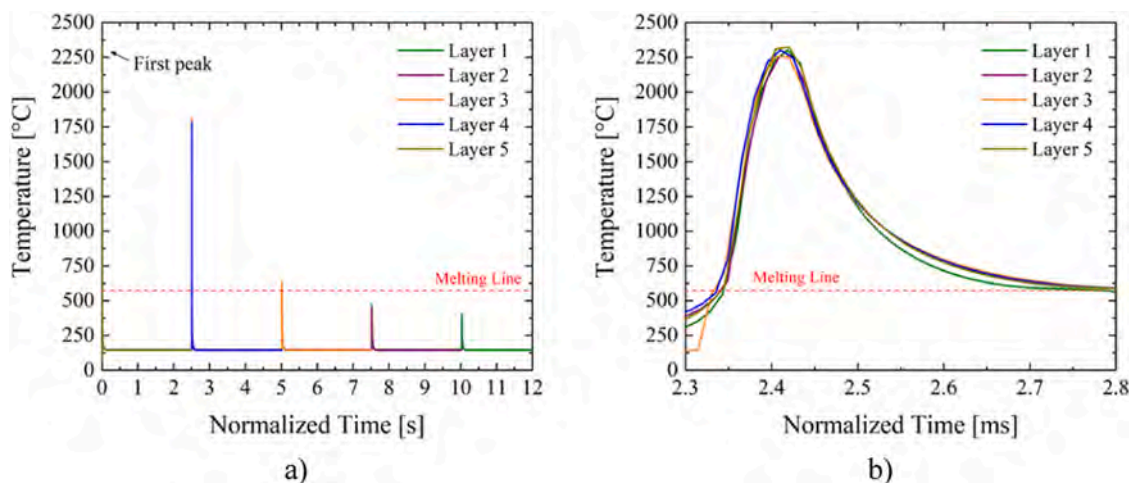


Fig. 15. Temperature cycle of different layers, (a) normalized thermal histories, (b) detail.

4. Conclusions

In this work, a 3D transient computational model was developed in the commercial FEM code MSC Marc to investigate the temperature field evolution during a LPBF process on AlSi10Mg alloy. The numerical results of a multi-track and multi-layer simulation were discussed regarding the experimental evidence. To validate the computational model, the predicted melt pool size was compared to the experimental one, and the results were satisfactory, with an error in the prediction of both the width and the depth less than 2%. OM and SEM analysis underlined a strong dependence of AlSi10Mg microstructure on the manufacturing process. The macrostructure shows a distinctive layered texture with overlapping melt pool profiles that represent the scanning path followed by the laser heat input. The major microstructural heterogeneities were found within the melt pool profile, with a variable cell size between the core and the periphery. Those features were linked to the thermal evolutions predicted by the model. Indeed, the temperature profiles of different points in a multi-track or a multi-layer analysis were comparable. Meanwhile, dissimilar thermal evolutions were experienced at different points of the melt pool, where the major thermal gradients can be found.

Declaration of Competing Interest

The authors declare that they have no known competing financial interests or personal relationships that could have appeared to influence the work reported in this paper.

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