

DUAL-OBJECTIVE OPTIMIZATION OF PBF-LB PROCESS VIA ENERGY DENSITY CONTROL

This study introduces a dual-objective optimization framework for the laser powder bed fusion (PBF-LB) of TC11 titanium alloy, designed to concurrently mitigate deformation and improve densification. A multiphysics finite element model, which incorporates temperature-dependent material properties, was developed to investigate the deformation mechanisms across an energy density range of 21.2-54.5 J/mm³. Experimental validation, conducted using 3D laser scanning, demonstrated close alignment with the predicted outcomes, with maximum deviations ranging from 0.04 to 0.08 mm. Systematic density measurements uncovered the nonlinear coupling effects between laser power and scan speed, resulting in an empirical power-velocity equation ($P = 0.2V + 30$, where P denotes laser power and V represents scanning speed), which yielded a peak relative density of 99.2%. The proposed strategy offers a comprehensive approach to the precision manufacturing of complex PBF-LB components, effectively balancing dimensional accuracy with mechanical performance, while minimizing experimental effort.

Keywords: Laser powder bed fusion; TC11 titanium alloy; Process parameter optimization; Deformation; Density

1. Introduction

Laser powder bed fusion (PBF-LB) is a near-net-shape additive manufacturing process widely adopted in the aerospace sector for producing high-performance titanium alloy components such as impellers and turbine blades. It offers short production cycles, high material utilisation, and the capability to fabricate complex geometries [1,2]. However, the nonlinear coupling of process parameters – laser power (P), scanning velocity (V), layer thickness (h), and hatch spacing (d) – creates dual challenges: controlling macroscopic deformation and achieving microstructural densification. Excessive energy input causes thermal stress accumulation (dimensional deviation >0.5 mm), while insufficient energy density results in unfused porosity (>5%) [3,4].

In addition to these parameters, scanning strategy [5], substrate preheating [6], atmosphere composition [7], and powder characteristics [8] also affect part quality. Although conventional optimisation approaches, such as orthogonal experiments [9,10] and response surface methodology [11,12], can reduce testing costs, they often fail to balance dimensional accuracy and mechanical performance, limiting PBF-LB applications in precision dynamic components [13].

Volumetric energy density ($E = P/(V \cdot h \cdot d)$) is frequently used as an integrated parameter for process optimisation and

cross-study comparison [14,15]. However, studies have shown that even at constant E , variations in power and speed combinations significantly alter microstructure and properties in Ti-6Al-4V and Al-12Si alloys [16,17], indicating that E alone cannot reliably predict material performance.

To overcome these challenges, this study introduces a dual-objective optimisation framework for PBF-LB of TC11 titanium alloy impellers, featuring two core innovations: (1) quantitative assessment of energy density (21.2-54.5 J/mm³) effects on impeller deformation via thermomechanical modelling, enabling accurate prediction; and (2) establishment of an empirical power-velocity relation ($P = 0.2V + 30$) for melt pool stabilisation, achieving 99.2% relative density in thin-walled regions (≤ 1 mm) and surpassing conventional E -based models.

2. Materials and methodology**2.1. Materials**

The TC11 titanium alloy powder used in this study was supplied by Xi'an OZ Technology Co., Ltd. and produced via the plasma rotating electrode process (PREP) [18]. It is an $\alpha + \beta$ dual-phase titanium alloy with a nominal composition of

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Ti-6.5Al-3.5Mo-1.5Zr-0.3Si (wt.%). The detailed chemical composition and particle size distribution data are provided in Tables S1 and S2 of the Supplementary Material. The morphology of the TC11 powder is shown in Figs. S1 to S3 of the Supplementary Material. The powder exhibited a particle size range suitable for laser powder bed fusion (PBF-LB), with characteristic particle diameters conforming to standard specifications for the process. Figs. S4 and S5 of the Supplementary Material show the experimental equipment and sample preparation platform, while Fig. S6 presents the final fabricated sample, the PBF-LB formed TC11 titanium alloy impeller.

2.2. Multiphysics modeling workflow

A comprehensive multiscale modeling framework was developed to simulate the macro-micro coupled behavior during the laser powder bed fusion (PBF-LB) of TC11 titanium alloy. Temperature-dependent thermophysical properties, spanning from 25°C to 1600°C (at 100°C intervals), were computed using JMatPro 11.0 [19], which has been validated for predicting thermal conductivity in titanium alloys with an error margin of less than 5% [20]. The obtained properties included thermal conductivity (5.8-24.6 W/m·K) [21], specific heat capacity (520-830 J/kg·K), and the coefficient of thermal expansion (9.2×10^{-6} - $11.7 \times 10^{-6}/^{\circ}\text{C}$), as shown in Fig. 1. The β -transus temperature was calculated to be 998°C through thermodynamic

analysis, and the latent heat of phase transformation (419 kJ/kg) was incorporated for subsequent thermal stress analysis.

The computed parameters were integrated into the Simufact Additive 2022 material database to develop a TC11-specific material model. The impeller geometry was discretized using a voxel-based mesh (voxel size = 1 mm; refer to Fig. 2), resulting in approximately 2.1×10^6 hexahedral elements. The volumetric energy density is defined as:

$$E = \frac{P}{V * d * h} \quad (1)$$

Where E represents the energy density (J/mm^3), P denotes the laser power (W), V is the scan speed (mm/s), d is the hatch spacing (mm), and h is the layer thickness (mm). The hatch spacing and layer thickness were fixed at 0.11 mm and 0.06 mm, respectively, ensuring consistency in the experimental setup. The laser spot diameter was set between 70 and 100 μm , and the substrate preheating temperature was maintained at 80°C. A Python automation script was employed to generate parameter sets spanning energy density levels from 21.2 to 54.5 J/mm^3 , requiring a total computational time of 187 hours.

2.3. Experimental design

A two-stage experimental protocol was developed to validate the multiphysics model within the dual-objective optimiza-

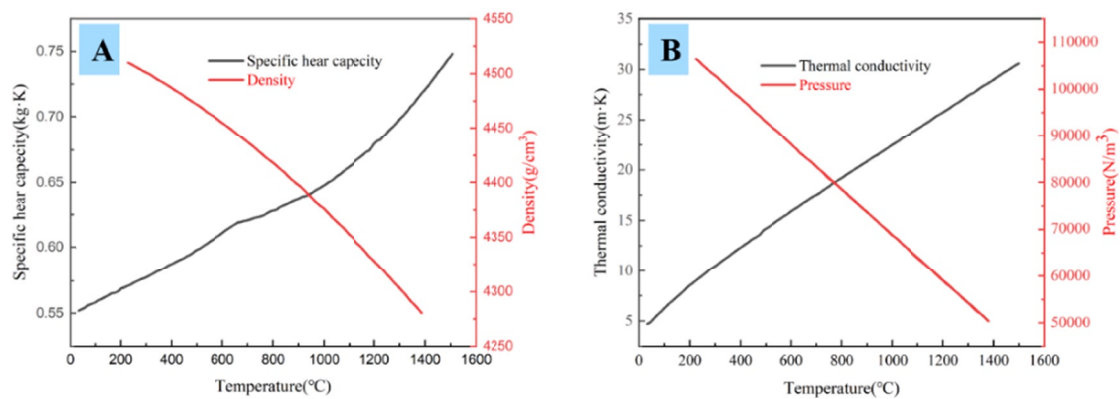


Fig. 1. JMatPro Thermophysical Properties of Titanium Alloy TC11: A – Density and Specific Heat Capacity vs. Temperature Curves; B – Thermal Conductivity and Young’s Modulus vs. Temperature Curves

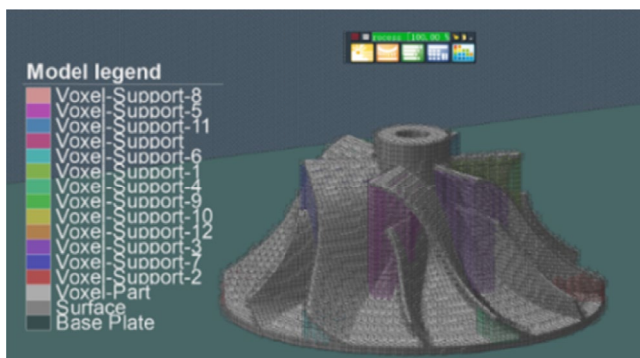


Fig. 2. Simufact model Voxel delineation map

tion framework. Stage 1: Energy Density-Deformation Mapping. This stage aimed to establish a quantitative relationship between volumetric energy density and the resulting deformation. Six energy density levels (21.2-54.5 J/mm^3) were systematically generated by varying laser power (210-360 W) and scan speed (1000-1500 mm/s), while layer thickness ($h = 0.06$ mm) and hatch spacing ($d = 0.11$ mm) were kept constant (see TABLE 1). Each parameter set was tested in triplicate to ensure statistical reliability and consistency. After printing, supports were removed via wire EDM, and full-field deformation measurements were captured using 3D laser scanning (Keyence VR-3200, accuracy ± 5 μm) across critical regions, including blade tips and

overhangs. The relative error between simulated and measured dimensions was calculated as:

$$\theta = \pm \frac{N-M}{N} \times 100\% \quad (2)$$

where M denotes the mean measured dimension and N the nominal dimension.

Stage 2: Densification-Oriented Optimization. This stage was focused exclusively on maximizing relative density through a structured parameter optimization approach. A full factorial design was employed, featuring six levels of laser power (210-360 W) and scan speed (1000-1600 mm/s), resulting in 36 distinct parameter combinations (TABLE 2). Cubic specimens ($15 \times 15 \times 10 \text{ mm}^3$) were fabricated in randomized positions on the build platform to minimize any bias resulting from localized thermal history. Density was measured using Archimedes' principle (ASTM B962), with five replicates per parameter set to ensure statistical robustness.

Validation trials on critical parameters ($P = 300 \text{ W}$, $V = 1200 \text{ mm/s}$, $E = 37.88 \text{ J/mm}^3$) were conducted using ANOVA ($\alpha = 0.05$) to confirm statistical consistency between simulated and experimental data, with error bands defined by the 3σ criterion (99.7% confidence).

TABLE 1

Experimental process parameters

Experiment number	Laser power (W)	Scanning speed (mm/s)	Energy density (J/mm^3)
1	210	1500	21.212
2	240	1400	25.974
3	270	1300	31.469
4	300	1200	37.879
5	330	1100	45.455
6	360	1000	54.545

TABLE 2

Optimized parameters for PBF-LB forming process of TC11 titanium alloy

Process parameters	Parameters
Laser power P (W)	250, 270, 290, 310, 330, 350
Scanning speeds V (mm/s)	1100, 1200, 1300, 1400, 1500, 1600
Powder thickness h (mm)	0.02, 0.04, 0.06
Scanning distance d (mm)	0.09, 0.11, 0.13

3. Results and validation

3.1. Validation of energy density-deformation relationship

Energy density gradient experiments revealed nonlinear deformation evolution in TC11 impellers (Fig. 3). When energy density increased from 21.2 J/mm^3 to 54.5 J/mm^3 , the maximum tip deformation (δ_{\max}) escalated from 0.29 mm to 0.73 mm , following the exponential model:

$$\delta_{\max} = 0.021e^{0.038E} (R^2 = 0.983)$$

3D laser scanning validation (Fig. 4) demonstrated a root mean square error (RMSE) of 0.059 mm between simulated and experimental data. The maximum absolute deviation occurred at $E = 25.9 \text{ J/mm}^3$ ($\Delta = 0.08 \text{ mm}$), corresponding to a relative error of 26.7% .

ANOVA analysis ($\alpha = 0.05$, $F = 2.34$, $p = 0.12$) confirmed no statistically significant difference between simulated and experimental data, with all error points residing within the 3σ confidence bands ($\pm 0.129 \text{ mm}$), thereby validating the model reliability.

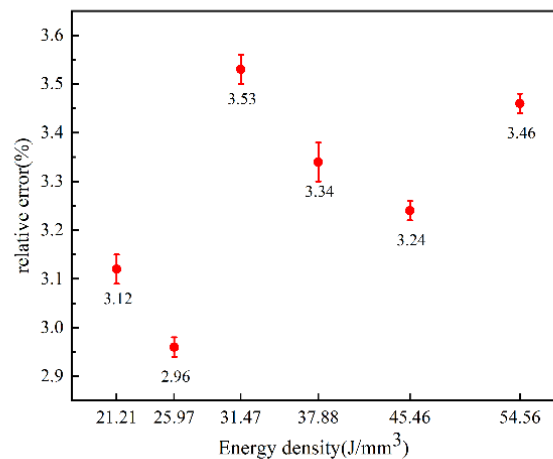


Fig. 3. Relative blade error for each energy density

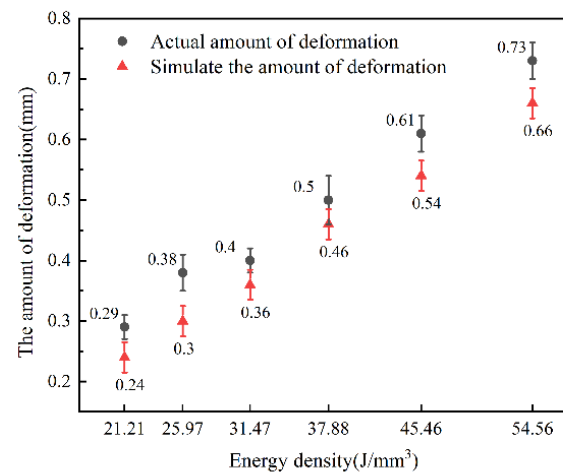


Fig. 4. Plot of energy density versus maximum simulated and actual deformations

3.2. Mechanism of process parameters on densification

The influence of process parameters on the relative density of TC11 titanium alloy during laser powder bed fusion (PBF-LB) is primarily governed by the dynamic regulation of melt pool behavior, which is modulated by energy density (E) [21]. Melt pool stability, which exhibits a high sensitivity to energy input, plays a critical role in dominating defect formation mechanisms. Particularly in keyhole mode (typically at $E > 50 \text{ J/mm}^3$), evap-

oration-induced recoil pressure can destabilize the melt pool, thereby promoting pore formation [22]. Experimental results demonstrate a nonlinear relationship between energy density and relative density. As energy density increases from 21.21 J/mm^3 to 54.5 J/mm^3 , the relative density improves nonlinearly from 96.852% to 99.197%, thereby confirming the existence of a critical energy density threshold (approximately 35 J/mm^3) required for adequate inter-layer bonding. Fig. 5 illustrates that at a fixed scan speed ($V = 1400 \text{ mm/s}$), relative density initially increases and then decreases with increasing laser power (P), peaking at $P = 310 \text{ W}$ (density = 99.197%). Fig. 6 further illustrates that when $P = 310 \text{ W}$, the maximum density is attained at $V = 1400 \text{ mm/s}$, validating the synergistic match between power and speed for optimal energy coupling.

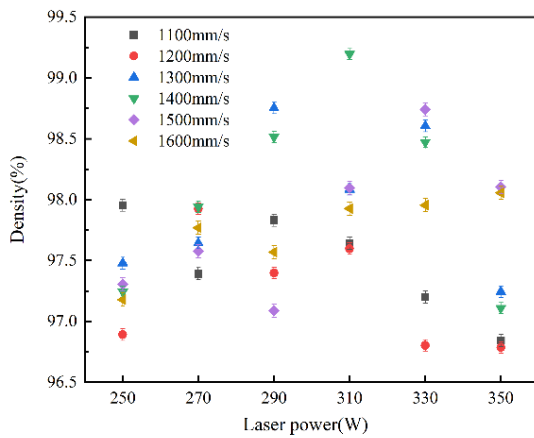


Fig. 5. Plot of laser power P versus densities

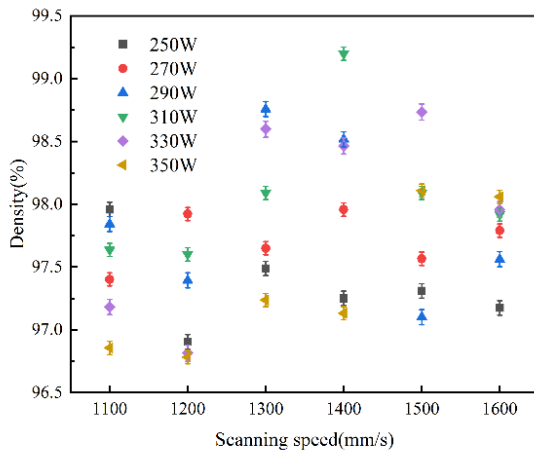


Fig. 6. Plot of scanning speed V versus densification

In the low-energy regime ($E < 35 \text{ J/mm}^3$), discontinuous melt pools result in unmelted defects, thereby reducing density below 98.5%. Within the optimized energy range ($35\text{--}42 \text{ J/mm}^3$), stabilized melt pools (aspect ratio ≈ 0.8) effectively suppress pore formation. At $E > 45 \text{ J/mm}^3$, melt turbulence induces the resurgence of gas pores, resulting in a significant decline in density. Parameter sensitivity analysis reveals that scan speed (V) exerts a dominant effect on density relative to laser power (P) and layer thickness (h), while P-h interactions exacerbate pore heteroge-

neity. Fig. 7 demonstrates that a hatch spacing of $d = 0.11 \text{ mm}$ (overlap ratio $\approx 30\%$) results in the highest density (99.1%). Deviations ($d > 0.12 \text{ mm}$ or $d < 0.10 \text{ mm}$) reduce density by more than 2.5%, primarily due to lack-of-fusion or remelting defects.

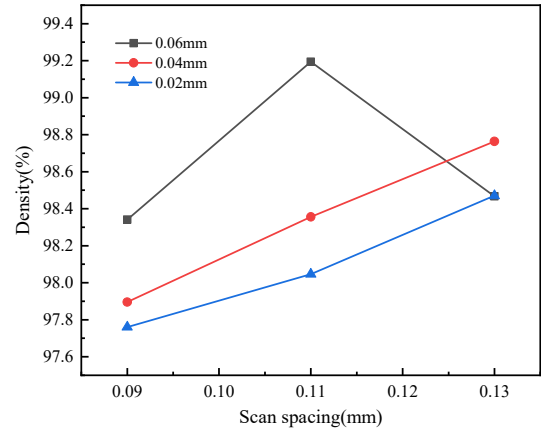


Fig. 7. Plot of scanning spacing d versus densification

The recommended process window includes: $P = 300\text{--}350 \text{ W}$, $V = 800\text{--}1200 \text{ mm/s}$, and $h = 30\text{--}50 \mu\text{m}$ ($V < 800 \text{ mm/s}$), supplemented by 67° interlayer rotation [23], which reduces anisotropic residual stress by 30% compared to unidirectional scanning [24], along with 200°C substrate preheating to mitigate keyhole instability. This strategy establishes a robust framework for the high-density PBF-LB fabrication of TC11 alloy components.

4. Conclusions

This study proposes a cost-effective process optimization framework for laser powder bed fusion (PBF-LB) of TC11 titanium alloy impellers, facilitating the synergistic control of both deformation and densification. The key findings are summarized as follows:

Through finite element simulations and 3D laser scanning experiments, a quantitative relationship between energy density ($E = P/(V \cdot d \cdot h)$) and residual deformation was established, with the maximum predicted deviation not exceeding 0.08 mm. At identical energy densities, deformation variations resulting from fluctuations in a single parameter remained below 3%, demonstrating that energy density serves as a robust and unified indicator for deformation control.

Laser power and scanning speed exert a significant influence on densification. Based on the empirical formula for melt pool stability regulation, $P = 0.2V + 30$, a relative density of 99.2% was achieved at $P = 310 \text{ W}$ and $V = 1400 \text{ mm/s}$, reducing process debugging efforts by 60% compared to traditional trial-and-error methods.

Within the energy density range of $35\text{--}42 \text{ J/mm}^3$, relative densities exceeding 99.0% and deformations below 0.15 mm were attained, establishing a viable process window for the direct fabrication of precision components, such as aviation engine combustion chamber liners.

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