



# Experimental investigation of the convective heat transfer coefficient for open-cell porous metal fins at low Reynolds numbers



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## ABSTRACT

The heat transfer characteristics and convective heat transfer coefficient of porous metal fin are experimentally investigated. The pore density of the porous metal fin and frontal velocity of the working fluid are varied in the range of 20, 40, and 80 pores per inch, and 0.007–0.17 m/s, respectively. The porous metal fins are fabricated from nickel with different porosities and various pore densities. The geometrical parameters of the test samples are measured using an optical method. In this study, porous metals are considered as a fin and the heat transfer performances are experimentally evaluated. An equivalent diameter based on the permeability and porosity is used as the characteristic length to calculate the Reynolds numbers and Nusselt numbers. When the equivalent diameter is used as the characteristic length, the measured Nusselt numbers converge to a single curve regardless of the pore density. The Nusselt number variation appears to be very similar to previous correlations developed for convective heat transfers of turbulent pipe flows. Consequently, an empirical correlation of the Nusselt number for porous metal fins is proposed in the form of a Dittus–Boelter correlation.

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## 1. Introduction

A heat exchange between working fluids is indispensable in various industrial and domestic energy systems. Because the heat exchange processes affect the overall performance of a thermal system, significant research results regarding compact heat exchangers have been presented over the past decade [1–4]. Porous metals have also garnered research interest for applications in electronic device cooling, fuel cell electrodes [5], compact heat exchangers [6], and so on. The diverse research areas in porous metals have been being investigated for more than 150 years [7–9]. Porous metals have high surface area-to-volume ratios as well as complex path structures. It is expected that enhancing the heat transfer area and flow mixing would increase the convective heat transfer capacity when applied to heat exchangers. Therefore, the heat transfer characteristics of porous metals have been investigated [10].

Rohsenow and Hartnett [11] and Nield and Bejan [12] considered forced convective heat transfers in rectangular ducts packed with porous materials and defined a convective heat transfer coefficient and Nusselt number based on the channel wall area. In their formulation, the overall porous medium thermal conductivity, i.e.

the weighted mean of the fluid thermal conductivity and solid thermal conductivity, was used to consider the effect the porous medium on the heat transfer augmentation. Most previous research concerning porous metals has used the local thermal equilibrium (LTE) model [13–15]. The LTE model assumes a local thermal equilibrium between the solid and fluid. Dukhan et al. [16] presented a macroscopic lumped-parameter model that obtains the temperature distribution in a rectangular block of thin porous metal foam with a constant cross-sectional area. However, the validity of the LTE model is restricted to low Reynolds number applications [17]. Whitaker [18] and Quintard and Whitaker [19] presented the constraints that must be satisfied in order for the LTE assumption to be valid. They demonstrated that the LTE model is only applicable when the conductive heat transfer is dominant in a representative elementary volume that consists of a fluid and a solid of porous metal. However, in most practical applications, the heat transfer to the fluid flowing through porous metals is dominated by convective heat transfers.

When the temperature between the solid and fluid in the porous metal differs, the LTE assumption cannot be applied [20]. The local thermal non-equilibrium (LTNE) model can be used when the temperatures of the fluid and porous metals are known and the heat transfer area is measured. For these reasons, many researchers have used the LTE model without verifying the adequacy of the LTE assumptions. The LTNE model was recently

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