



Theoretical and experimental studies on a solid containing water droplet



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ARTICLE INFO

Article history:

Received 1 January 2014

Received in revised form 20 June 2014

Accepted 22 June 2014

Available online 12 July 2014

Keywords:

Dissolved solids

Single droplet

Water evaporation

Glass filament method

Spray cooling

ABSTRACT

Heat and mass transfer to and from a single solution droplet is studied in this work. A new theoretical model to predict the evaporation behaviour of solid containing water droplets is presented. The model, implemented in MATLAB, is used to predict the process of droplet evaporation with prediction results successfully validated against data from the literature. Also, an experimental study was performed to study the evaporation of a single droplet containing NaCl and water. To investigate the influence of concentration, tests were performed with droplet having initial radiuses of approximately 0.5 mm and initial mass concentrations of 3% and 5%. Results obtained from the developed model were found to be in good agreement with our experimental data. Finally, it was shown that the current model, allowing for a smooth transition from surface evaporation to crystallisation, is able to simulate the process more accurately compared to existing models in the literature which lead to a, less realistic, sharp transition.

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1. Introduction

Liquid sprays are widely applied in engineering applications with cooling towers, evaporators, condensers, and chemical reactors as some examples. The benefit of using these liquid sprays is increasing the contact area to obtain higher total heat and mass transfer between liquid and air. Applying this technique in arid areas, however, faces an obvious challenge of providing pure water. Hence, saline water is being used as an alternative in the spray-assisted dry (hybrid) cooling towers. Heat exchanger surfaces are in exposure of corrosion and deposit formation due to using saline water because of large amount of corrosive ions in water. Therefore, contact of solid containing water droplets in a hybrid cooling tower to the heat exchanger surfaces should be avoided. In other words, evaporation of the droplets should be completed before they reach the heat exchanger surfaces.

Beginning to evaporate, the evaporation rate of a solid-containing droplet is the same as that of pure water droplet. However, when the average concentration (mass fraction of the solid divided by the total mass of the droplet) reaches that of saturation, solid formation occurs leading to an increase in the droplet temperature [1]. This temperature rise is due to both heat of crystallisation and sensible heat transfer in the case of solutions, and to sensible heat transfer in the case of suspensions [1].

1.1. Review of existing evaporation models

1.1.1. Two-stage models

There are models in the literature that assume two stages for droplet evaporation. According to those models, during the first stage of evaporation the main component of the droplet is assumed to be water. The second stage begins when the solid content of the droplet increases, due to water evaporation, and goes beyond a critical concentration value [2].

Applying a two-stage model, Abuaf and Staub simulated the latter stage by assuming the droplet as a porous solid crust around a wet core at the centre [2]. Elperin and Krasovtsov assumed quasi-steady evaporation and showed that the evaporating rate is highly dependent on the permeability of the formed porous crust for vapour and heat transfer [3].

1.1.2. Three-stage models

On top of two-stage models, some researchers showed that three different stages occur during evaporation of a solution droplet [3–6]. According to those authors, the first stage of the previous model is divided into two stages including: temperature adjustment stage and constant-temperature evaporation stage.

Three-stage models assume that the droplet temperature remains constant beyond the wet-bulb temperature. This isothermal stage is then followed by a sharp rise in the droplet temperature to reach the ambient air temperature. Also, a d^2 law is assumed for the time taken for the onset of crust formation [7]. A mathematical model was presented by Farid to predict the

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