

CFD Analysis of nanofluid forced convection heat transport in laminar flow through a compact pipe

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Abstract. In the present paper, developing laminar forced convection flows were numerically investigated by using water-Al₂O₃ nano-fluid through a circular compact pipe which has 4.5mm diameter. Each model has a steady state and uniform heat flux (UHF) at the wall. The whole numerical experiments were processed under the $Re = 1050$ and the nano-fluid models were made by the Alumina volume fraction. A single-phase fluid models were defined through nano-fluid physical and thermal properties calculations, Two-phase model(mixture granular model) were processed in 100nm diameter. The results show that Nusselt number and heat transfer rate are improved as the Al₂O₃ volume fraction increased. All of the numerical flow simulations are processed by the FLUENT. The results show the increment of thermal transfer from the volume fraction concentration.

1. Introduction

To enhance the heat exchangers energy transfer efficiency, Engineers have interested not only the development exchanger or enhanced its structure, but also improvement of heat transfer fluid itself. Maxwell [1] proved the possibility of increasing thermal conductivity of a mixture by higher volume fraction of solid particle. Nowadays available particle diameter is smaller than 100nm which is called 'Nano-sized'. This effort makes particle mixture fluid would enhance the heat transfer performance of heat transfer fluids. [2] Moreover, solid nanoparticle colloids are highly stable and exhibit no significant settling under static conditions. Also, decreasing particle size makes suspending solid colloid easily. So that nano-fluid technology is expected to suitable of heat transfer fluid. [3] However, nano-fluid development is interrupted by the lack of agreement with different research group, theoretical understanding of the heat transfer mechanisms, different suspension conditions. [4] In this paper, Nano-fluid model assume that the Homogeneous flow mixture model [5] which is the dispersed and continuous phase with very strong coupling and the nano-particles is moving at the same inlet velocity. Therefore instead of DPM, the dispersed granular mixture models are made by 'Mixture model' which is one of the Euler-Euler approaches. This numerical investigation purpose is an observation of heat transfer coefficient increment in a compact pipe and comparison between single-phase and multi-phase model. Geometry model size is as same as experimental pipe which will use in



a university laboratory. Nano-particle volume fractions are less than 5 percent, and divided by 1 percent from 0 to 4 percent. The wall has constant surface heat flux as known as Uniform Heat Flux (UHF). The numerical simulation geometry is axisymmetric two-dimensional model. Through the limit of 2D model and assumption, implicit effect such as gravity and buoyancy force are excluded. The nano particle material is an aluminum oxide with a spherical size of 100 nm diameter and simulations are processed in a steady state.

2. Mathematical modelling

2.1. Geometry and boundary condition

Fig. 1a shows the actual model of geometry and Fig. 1b is the part of the Two-dimensional Mesh. One of the main purpose of this simulation is trying to understand the near wall behavior of nano-particles, velocity distribution perpendicular to the pipe surface and drag coefficient near at the pipe wall, so the relatively concentrated mesh in the near wall area is needed. As a result, the mesh has a non-uniform quadratic mesh. The bottom horizontal line is an axis which makes Three-dimensional pipe model at the axisymmetric environment. The upper horizontal line is a wall with UHF. The tube consisted with a diameter (D) of 4.5 mm and a length (L) of 1.17 m. The nano-fluid is composed of water and Al_2O_3 particles. Whole velocity inlet for each fluid models enter with uniform temperature and axial velocity profiles are defined by $\text{Re} = 1050$. The fluid is flow out to the outflow outlet.

2.2. Governing equation

The single-phase model is investigated for the comparison of mixture model as prescribed. Following equations are representing the mathematical formulation of the single-phase model and of the continuous phase of the mixture model. [5-7]

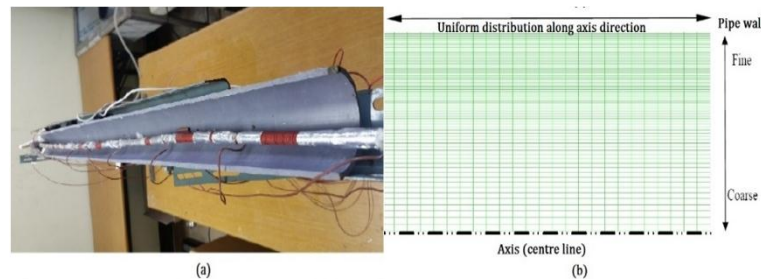


Figure 1. (a) Actual model of geometry (b) Part of the mesh of Computational Cell. 80 x 8000 with 0.96 ratio

Energy equation

$$\text{div}(\rho \vec{V} C_p T) = \text{div}(k \text{grad} T) + S_e \quad (1)$$

Conservation of mass

$$\frac{\delta \rho_m}{\delta t} + \nabla \cdot (\rho_m u_m) = 0 \quad (2)$$

Mass-averaged velocity u_m is

$$u_m = \frac{\sum_{k=1}^n \alpha_k \rho_k u_k}{\rho_m} \quad (3)$$

and ρ_m is the mixture density with α_k volume fraction of phase k

$$\rho_m = \sum_{k=1}^n \alpha_k \rho_k \quad (4)$$

Momentum equation for single-phase

$$\text{div}(\rho \vec{V} \vec{V}) = -\text{grad} P + \nabla \cdot (\mu \nabla \vec{V}) + S_m \quad (5)$$

Momentum equation for multi-phase

$$\nabla \cdot \sum_{k=1}^n \alpha_k \rho_k u_k u_k = \nabla \cdot (\rho_m u_m u_m) + \nabla \cdot \sum_{k=1}^n \alpha_k \rho_k u_{Mk} u_{Mk} \quad (6)$$

The compression and viscous lose by heat are assumed to be excluded in the energy equation. The source/sink terms S_m and S_e represent the unitive effects of energy and momentum exchange with base fluid. These are zero value in the single-phase model. Momentum equation of multi-phase are affected by volume fraction (α), velocity of phase (u_k), diffusion velocity (u_{Mk}). The determinations of particle

motion are particle density(ρ_c), cross-sectional area of the particle(A_p), drag coefficient(C_D), slip velocity(u_{cp}). The drag force in this study is following Clift et al equation (7) for single rigid spherical particle in a fluid

$$F_D = -\frac{1}{2}A_p\rho_c C_D |u_{cp}|u_{cp} - \frac{1}{2}V_p\rho_c \frac{du_{cp}}{dt} - 6\gamma_p^2 \sqrt{\pi\rho_c\mu_c} \int_0^t \frac{du_{cp}}{\sqrt{t-s}} ds \quad (7)$$

The drag coefficient which is already showed in equation (7) depends on small Particle Reynolds numbers with particle Reynolds number

$$C_{D,st} = \frac{24}{Re_p} \quad (8)$$

The slip velocity is following equation (9) with τ_p the particle relaxation time, \vec{a} is the secondary-phase particle's acceleration

$$u_{pq} = \frac{\tau_p(\rho_p - \rho_m)}{f_{drag}} \vec{a} \quad (9)$$

The default drag function is taken Schiller and Naumann. Reynolds numbers in this paper are over 1000 so the equation is defined as follow

$$f_{drag} = 0.0183Re \quad (10)$$

2.3. Physical properties of the nanofluid for single-phase model and Boundary conditions

At inlet boundary condition, axial velocity V_0 is defined from $Re = 1050$ according to the fluid properties and a constant temperature (293K) as described. On the tube wall, the usual non-slip conditions and uniform wall heat flux $q = 5000 \text{ W/m}^2$ are implemented.

The determination of nano-fluid physical, thermal properties are at the center of current nano-fluids research as previously mentioned.[7] For the single phase model properties of the nano-fluids have been calculated by using classical formulas developed for conventional solid-liquid mixtures.[8] The following equations are considered to evaluate Al_2O_3 -water nano-fluids thermo-physical properties and were used for CFD simulation.[6]

$$\rho_{nf} = (1 - \varphi)\rho_{bf} + \varphi\rho_p \quad (11)$$

$$c_{p,nf} = \frac{(1 - \varphi)(\rho c_p)_{bf} + \varphi(\rho c_p)_p}{\rho_{nf}} \quad (12)$$

$$\mu_{nf} = (123\varphi^2 + 7.3\varphi + 1)\mu_{bf} \quad (13)$$

The φ is the volume fraction of the Al_2O_3 . The dynamic viscosity that temperature depended properties are excluded and has a constant value. The equation (13) is applicable for spherical particle in volume fractions less than 5% [9] and determination of nano-fluid effective thermal conductivity as follow. [10]

$$k_{nf} = \left[\frac{k_s + 2k_w + 2(k_s - k_w)(1 + \beta)^3\varphi}{k_s + 2k_w - (k_s - k_w)(1 + \beta)^3\varphi} \right] k_w \quad (14)$$

In Equation (14), β is the ratio of nano-layer thickness to the original particle radius, and $\beta = 0.1$ is used for calculation in this equation. The result of physical and thermal properties of nano-fluid models are presented in Table 1 and Al_2O_3 particle physical and thermal properties as follow in Table 2

2.4. Numerical approach

The computational fluid dynamic code FLUENT 16.1 is employed to solve the present problem. The governing Eqs. (1)-(3) are solved by control volume approach. For the convective and diffusive terms, a second order upwind method was used. Pressure and velocity were coupled using Semi Implicit Method for Pressure Linked

Table1. Properties of nano-fluid at 293K and $Re = 1050$

$\Phi(\%)$	Density (kg/m^3)	Viscosity (kg/ms)	Inlet velocity (m/s)	Heat conductivity (W/mK)	Specific heat (J/kgK)
0	998.2	9.98e-4	0.233	0.5970	4182.0
1	1027.018	1.023e-3	0.2324	0.6200	4053.21
2	1055.836	1.048e-3	0.2316	0.6436	3931.451
3	1084.654	1.073e-3	0.2308	0.6678	3816.162
4	1113.472	1.098e-3	0.2300	0.6926	3706.840

Table 2. Al_2O_3 nanoparticle at temperature of 293K

Chemical formula	Density (kg/m ³)	Heat conductivity (W/mK)	Specific heat (J/kgK)	Spherical diameter (nm)
Al ₂ O ₃	3880	36	773	100

Equations (SIMPLE) in single-phase. Unlike single-phase, multi-phase was approached by Coupled algorithm which offers an alternative to the density-based and pressure-based segregated algorithm with SIMPLE-type pressure-velocity coupling. The difference between SIMPLE and Coupled is a Convergence time [6] and Couple algorithm has converged faster than SIMPLE in multi-phase. The laminar flow has fully developed region that distribution of flow velocity maintains constant. In this study, Fully developed region appears after 0.307m according to equation (15). [12].

$$L_{laminar} = 0.065 \cdot Re \cdot D \quad (15)$$

The velocity distributions are showed in Figure 2. It is only 0.6737 percent less than outlet at an axial velocity at the 0.307m in Fig. 2a and all mixture models are following the Fully developed region in Fig. 2b. This means nano-fluid motion is fully developed at exit section model, when properties are not dependent on temperature, velocity profile and hydrodynamic entrance length are nearly identical at 0.307m for $\Phi = 0\%$, 1%, 2%, 3% and 4%. [7]

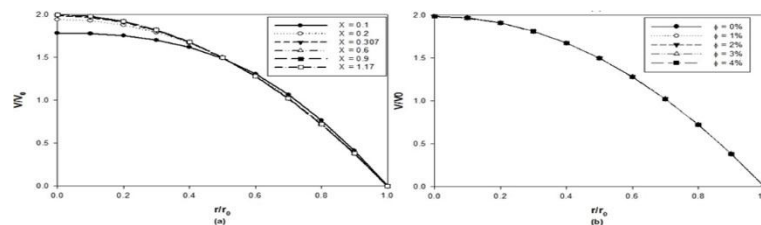


Figure 2. (a) Profiles of axial velocity at several locations for $Re = 1050$, $q = 5000 \text{ W/m}^2$, $\Phi = 4\%$. (b) Profiles of axial velocity at $x = 0.307\text{m}$ for $Re = 1050$, $q = 5000 \text{ W/m}^2$ and several volume fractions.

A Nusselt number is converged to 4.36 in the UHF.[13] Comparing Nusselt number to Churchill-Ozoe equation (16) [14] and Siegel Sparrow-Hallman calculation equation (17) [15] and common Nusselt number equation (18) which used in present study in Figure 3. x^* is a dimensionless axial position for thermally developing flows. Pr is the Prandtl number.

$$Nu = 5.364[1 + (70.0x^*)^{-10/9}]^{3/10} - 1 \quad (16)$$

$$Nu = \frac{2}{\left[\left(\frac{t_w - t_0}{qr_0/k}\right) - \left(\frac{4x/r_0}{RePr}\right)\right]} \quad (17)$$

$$Nu = \frac{h(x)D}{k_{inlet}} \quad (18)$$

The heat transfer coefficient $h(x)$ is defined as

$$h(x) = \frac{q}{T_w(x) - T_b(x)} \quad (19)$$

The $T(x)_w$ is the solid surface temperature at the wall temperature, and $T(x)_b$ is the mean bulk temperature. Average Nusselt number is defined as

$$Nu_{av} = \frac{h_{av} \cdot D}{k_{inlet}} \quad (20)$$

Current study model on $\phi = 0\%$ is going to 4.69 Nusselt number which has a 7 percentage point gap between normal convergence value. Therefore velocity distribution and Nusselt number show the model has fitted in the theory.

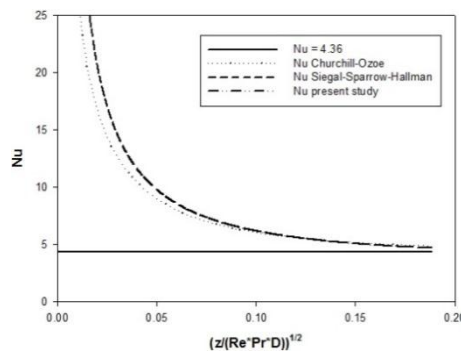


Figure 3. Grid validation by means of Churchill and Ozoe, Siegal-Sparrow-Hallman

3. Results

In this work a number of numerical simulations have been performed to study forced convective heat transfer of Al_2O_3 -water in a circular tube under laminar flow and results were conducted to employing the single phase and mixture model for $\phi = 1\%$ to 4% , $\text{Re} = 1050$ as $q = 5000 \text{ W/m}^2$ with a constant state. In all cases the size of the spherical particles is 100 nm. Dimensionless temperature shows the particle concentration increase happens the thermal development acceleration of the fluid at several axial locations along the radius is reported in Fig. 4 for $\phi = 0\%$ and 4% . For the both the base fluid and the nano-fluid, the motion is not thermally developed. As it is also noticed by a higher slope of relative local heat transfer coefficient h_r for $\phi = 4\%$ shown in Fig. 5a. Relative local heat transfer coefficient is increasing with axis location and volume fraction rise when fluid properties are constant in Fig. 5a. Because of narrow hydrodynamic radius of model, correlation between wall and bulk temperature is higher than larger model. Consequently, difference between the wall and bulk temperature is decrease and with the heat flux on the wall being constant, there is an increase in the heat transfer coefficient, as shown in Fig. 5b. Small displayed equations: Some examples:

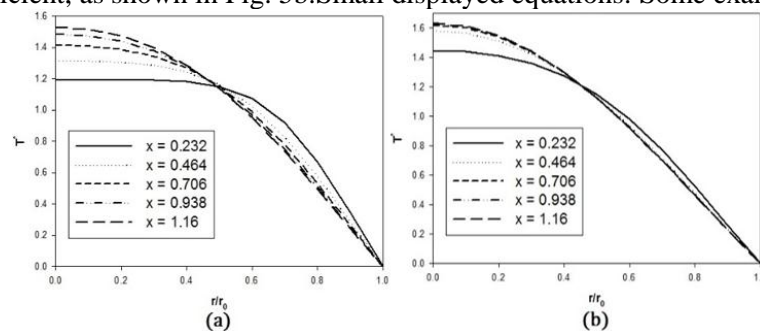


Figure 4. Dimensionless temperature for $\text{Re} = 1050$ and $q = 5000 \text{ W/m}^2$ at several locations and for : (a) $\Phi = 0\%$ and (b) $\Phi = 4\%$

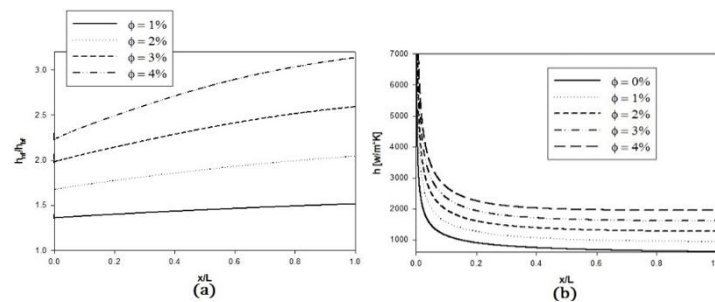


Figure 5. (a) Increase in nano-fluid heat transfer coefficient along tube axis for $\text{Re} = 1050$ and $q = 5000 \text{ W/m}^2$ for constant properties and increment of volume fraction. (b) Heat transfer coefficient of each volume fraction for constant properties.

In Fig. 6, radial temperature profiles at profiles at $x/L = 1.0$ for $\text{Re} = 1050$ and $q = 5000 \text{ W/m}^2$ are investigated for the effect of particle presence. The temperature is strongly dependent on ϕ

concentration. When the multi-phase concentration increases, the difference between temperature at surface and axis decrease by higher forced convection heat transfer. This effect is the cause of wall temperature decrease. It is also implemented in single-phase however, due to the particle absence, heat transfer between wall and inner side is not remarkable. Furthermore the only average temperature 1.3K increase than 293K on single-phase models. Both of phase show there is a clear existence of a uniform temperature fluid zone that becomes more visible for higher value of Φ at the $r/r_0 = 0.6$.

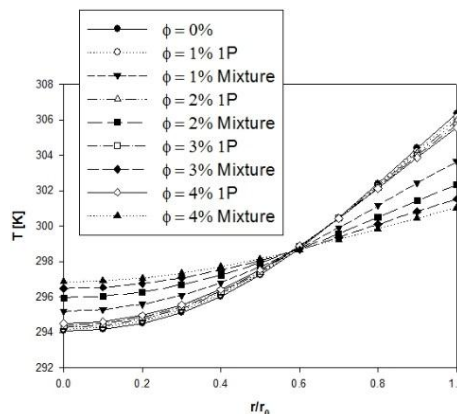


Figure 6. Comparison of temperature along tube radius at $x/L = 0.99$ for several concentration and for $Re = 1050$ and $q = 5000 \text{ W/m}^2$ in constant properties

Table3. Heat transfer coefficients for $Re = 1050$, $q = 5000 \text{ W/m}^2$ and the investigated models

$\Phi(\%)$	$h_{av} \text{ (W/mK)}$	h_{nf} / h_{bf}	Nu_{av}	Nu_{nf} / Nu_{bf}
0	861.815	1	6.5	1
1	1233.288	1.431	8.95	1.377
2	1592.853	1.848	11.1	1.707
3	1954.472	2.268	13.2	2.027
4	2301.521	2.671	14.9	2.302

4. Conclusion

Average heat transfer coefficient and Nusselt number are shown in Table 3 with relative increase of the total heat transfer rates and Nusselt number as a concentration of the nano-particle volume fraction ϕ . Significant increases of the total heat transfer rates can be found with the use of fully dispersed nanoparticles multi-phase. These results have indicated the nano-fluids beneficial effects of thermal properties improvement. The results showed that the nanoparticles inclusion produced considerable heat transfer with respect to base fluid. Heat transfer enhancement was increasing with concentration of particle volume. But the CFD should be progressed with constant and transient numerical investigation parallel and these numerical researches would be pace with the experiment especially. Also, lack of Reynolds number and uniform heat flux example, reality of nano-particle dispersed state are not including in this paper.

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