

Study of thermal conductivity of nanofluids for the application of heat transfer fluids

Dae-Hwang Yoo^a, K.S. Hong^b, Ho-Soon Yang^{c,*}

^a *Research Center for Dielectric and Advanced Matter Physics, Pusan National University, Busan 609-735, Republic of Korea*

^b *Busan Center, Korea Basic Science Institute, Busan 609-735, Republic of Korea*

^c *Department of Physics, Pusan National University, Busan 609-735, Republic of Korea*

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Abstract

TiO₂, Al₂O₃, Fe, and WO₃ nanofluids are prepared in a two-step procedure by dispersing nanoparticles in a basefluid. Since nanoparticles form clusters in fluids, a cell disrupter generating high power pulses is used for improving the dispersion of nanoparticles. The transient hot wire method is used for the measurement of thermal conductivity. The thermal conductivities of TiO₂, Al₂O₃, Fe, and WO₃ nanofluids are studied and compared with each other. Nanofluids show a large enhancement of thermal conductivity compared with their basefluids, which exceeds the theoretical expectation of two-component mixture system. We compare thermal conductivities of various nanofluids and discuss the important factors in determining thermal conductivity in this study.

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Keywords: Nanofluids; Thermal conductivity; Hot wire method; Ceramic nanoparticles; Metallic nanoparticles

1. Introduction

It has been recognized that the suspension of solids in fluids enhances the effective thermal conductivity of the material [1]. Enhancement of thermal conductivity of fluids contributes to improving the efficiency of heat transfer fluids. Furthermore, it is possible to reduce the size of the heat exchange system, which has been limited due to the poor thermal transport property of fluids. In earlier days, micron-sized or larger particles were suspended in fluids and they led to causing the problems, such as the settlement of particles, clogging, abrasion of devices, etc. [2].

It was proposed that fluids containing nanometer-sized particles can be a new class of engineered fluids with high thermal conductivity [2]. The nanoparticle suspended fluids, named nanofluids, have been produced as the nanotechnology producing nanoparticles developed rapidly. There have been many researches in the enhancement of thermal conductivity of nanofluids [3–8]. It was reported that the nanofluid containing 10-nm-sized Cu nanoparticles showed a large enhancement of thermal conductivity compared with Al₂O₃ nanofluid con-

taining 35 nm nanoparticles [6]. We also reported that 18% enhancement of thermal conductivity was achieved by using 10 nm Fe nanoparticles [7]. Large enhancement of thermal conductivity of nanofluids could not be predicted by conventional theories on the effective thermal conductivity of two-component materials [9–11]. The exact mechanism of thermal transport in nanofluids is not known at this moment, even if several potential mechanisms were suggested to describe experimental results of thermal conductivity of nanofluids. Many factors, such as particle size, effect of surfactant, dispersion of particles, and thermal property of dispersed particles have been expected to influence the thermal property of nanofluids. Moreover, the surface-to-volume ratio of particles increases dramatically as the particle size reduces, which can lead to potentiality of the significant enhancement of thermal conductivity of fluids [7]. The successful commercialization of nanofluids will bring the industrial advantages by reducing the energy consumption and the exhaust gas [12].

The results of Cu and Al₂O₃ nanofluids suggested the effect of intrinsic thermal conductivity and particle size of dispersed nanoparticles on thermal conductivity of nanofluids, but more nanofluids are needed for understanding of the effect. This work uses the oxide nanoparticles, such as TiO₂, Al₂O₃, and WO₃ and metal nanoparticles of Fe for nanofluids for the study of thermal conductivity of nanofluids. As comparing those results with

* Corresponding author. Tel.: +82 51 510 2221; fax: +82 51 515 2390.

E-mail address: hsyang@pusan.ac.kr (H.-S. Yang).

Al₂O₃, and Cu nanofluids, we would like to understand important factors in determining thermal conductivity of nanofluids. We also compare the experimental results with the calculated results with the conventional model for two-component materials.

2. Experimental

TiO₂, Al₂O₃, Fe, and WO₃ nanoparticles were synthesized by a chemical vapor deposition process for the preparation of nanofluids. The mean sizes of TiO₂, Al₂O₃, Fe, and WO₃ nanoparticles determined with transmission electron microscope (TEM) images are 25 nm, 48 nm, 10 nm, and 38 nm, respectively. The crystal structures of nanoparticles were determined with X-ray diffraction (XRD) patterns. We prepared TiO₂, Al₂O₃, Fe, and WO₃ nanofluids in a two-step procedure by dispersing nanoparticles in basefluids. Deionized water was used as a basefluid for TiO₂ and Al₂O₃ nanofluids, and ethylene glycol was used for Fe and WO₃ nanofluids. Table 1 shows the size and volume fraction of nanoparticles and basefluids of the nanofluids prepared in this study. Since no stabilizing agents were added in the preparation of nanofluids, nanoparticles must have been forming clusters in nanofluids. In order to improve the dispersion of particles in fluid, we used an ultrasonic cell disrupter (Jeiotech, ULH-700S) generating ultrasonic pulses of 700 W at 20 kHz. Even though the sonication cannot break nanoparticles individually, the cluster of nanoparticles breaks into smaller clusters resulting in the decrease of clusters [7,8].

Thermal conductivity of nanofluids was measured with a transient hot wire (THW) method, which is one of the most accurate methods for determining thermal conductivity of fluids [13,14]. A TeflonTM-coated pure (99.9%) platinum wire was used in the measurement. The wire is used as both heater and thermometer. THW method measures the temperature and time response of the wire to an abrupt electrical pulse. Thermal conductivity, k , is calculated from a derivation of Fourier's law:

$$k = \left[\frac{q}{4\pi(T_2 - T_1)} \right] \ln \left(\frac{t_2}{t_1} \right), \quad (1)$$

where q is the applied electric power and T_1 and T_2 are the temperatures at times t_1 and t_2 , respectively. From the temperature coefficient of the electrical resistance of a wire, the temperature rise of the wire can be determined by the change in its electrical resistance with time [12]. We measured thermal conductivity of nanofluids right after the sonication for 50 min. The thermal conductivity of nanofluids was obtained as a function of volume fraction of nanoparticles suspending in the nanofluids.

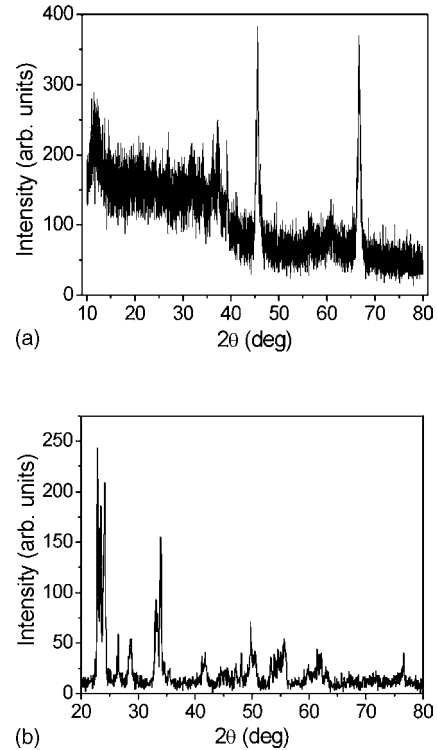


Fig. 1. XRD patterns of: (a) Al₂O₃ and (b) WO₃ nanoparticles.

3. Results and discussion

XRD patterns of Al₂O₃ and WO₃ nanoparticles are given in Fig. 1(a and b), respectively. Al₂O₃ nanoparticles form in γ -Al₂O₃ phase having the peaks of (4 0 0) and (4 4 0) and WO₃ nanoparticles form in monoclinic phase. The main XRD peaks at 23° and 24° correspond to the (0 0 2) and (1 1 0) reflections in bulk WO₃ crystal.

Fig. 2 shows the thermal conductivities of TiO₂ and Al₂O₃ nanofluids as a function of volume fraction of nanoparticles. Deionized water was used as a basefluid for both nanofluids. k and k_0 in the figure represent thermal conductivity of nanofluid and basefluid, respectively. As reported by other groups [6], Al₂O₃ shows very slow increase with the volume fraction of nanoparticles. It has been known that nanofluids containing ceramic nanoparticles exhibit smaller enhancement compared with nanofluids containing metallic nanoparticles. While there is only 4% enhancement of thermal conductivity with 1.0% volume fraction of Al₂O₃ nanoparticles, TiO₂ containing 1.0% volume fraction of nanoparticles exhibits the 14.4% enhancement of thermal conductivity. TiO₂ nanofluid shows higher

Table 1
Characterization of nanoparticles and basefluids used in the preparation of nanofluids

	Nanoparticle			
	TiO ₂	Al ₂ O ₃	Fe	WO ₃
Average size of particles (nm)	25	48	10	38
Volume fraction	0.1, 0.5, 1.0	0.3, 0.5, 0.7, 1.0	0.2, 0.3, 0.4, 0.55	0.05, 0.1, 0.2, 0.3
Basefluid	Water		Ethylene glycol	

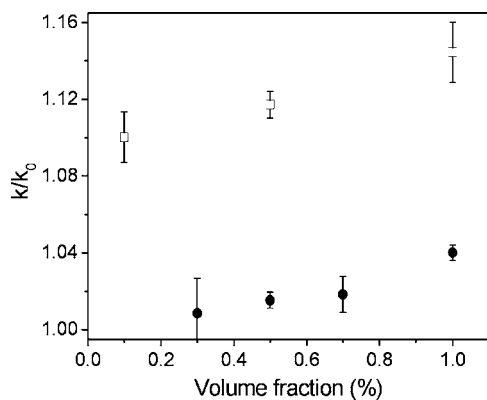


Fig. 2. Thermal conductivities TiO₂ (opened squares) and Al₂O₃ (closed circles) nanofluids as a function of volume fraction. Deionized water is used as a basefluid for both nanofluids.

enhancement of thermal conductivity compared with Al₂O₃ nanofluids even though bulk TiO₂ crystal shows lower thermal conductivity than bulk Al₂O₃ crystal. In our previous works for nanofluids, we found that the suspension of highly thermal conductive materials does not directly lead to the larger enhancement of thermal conductivity of nanofluids [8]. The comparison of Al₂O₃ and TiO₂ nanofluids supports that thermal conductivity of nanoparticles is not a principal factor related to thermal conductivity of nanofluids. Since the particle size of TiO₂ is smaller than that of Al₂O₃, the thermal conductivity of TiO₂ is larger than that of Al₂O₃ due to the larger surface-to-volume ratio.

Fig. 3 shows the thermal conductivities of Fe and WO₃ nanofluids as a function of nanoparticle volume fraction, where ethylene glycol was used as a basefluid. Fe and WO₃ nanofluids show the different trend in increase of thermal conductivity. Fe nanofluid shows 16.5% enhancement of thermal conductivity as containing 0.3% volume fraction of nanoparticles while WO₃ nanofluid of 0.3% volume fraction of nanoparticles shows 13.8% enhancement, which is consistent with other reports that ceramic nanoparticle-suspending nanofluids exhibited lower thermal conductivity than metallic nanoparticle suspending nanofluids. We also keep in mind that the size of Fe nanoparticles used in this study is smaller than that of WO₃. From the results of Figs. 2 and 3, we understand that the surface-to-volume ratio

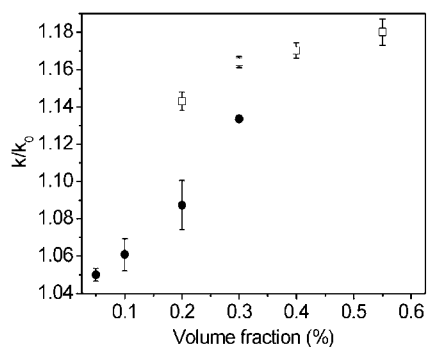


Fig. 3. Thermal conductivities of Fe (opened squares) and WO₃ (closed circles) nanofluids as a function of volume fraction. Ethylene glycol is used as a basefluid for both nanofluids.

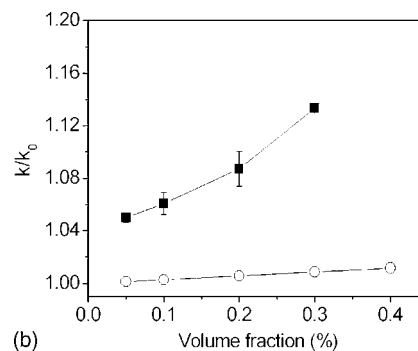
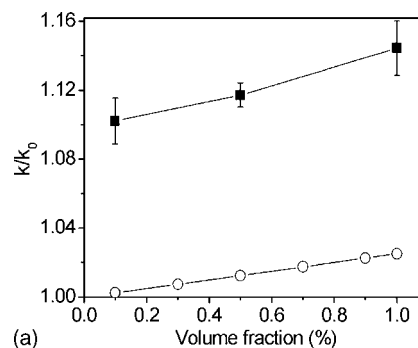


Fig. 4. The calculated thermal conductivities with the Hamilton–Crosser model for: (a) TiO₂ and (b) WO₃ nanofluids. Open circles represent the calculated values and closed squares represent the experimental results.

of nanoparticles is a principle factor in determining thermal conductivity of nanofluids.

For two-component mixtures with the micrometer-sized particles or larger particles, there are many models for thermal conductivity of the mixtures and they coincide with the experimental results [9–11]. We calculated thermal conductivity of TiO₂ and WO₃ nanofluids using Hamilton–Crosser (H–C) model which is one of the conventional models for two-component mixtures [10]. Water and ethylene glycol were chosen as a basefluid for TiO₂ and WO₃ nanofluids, respectively. Fig. 4(a and b) presents the thermal conductivities of TiO₂ and WO₃ nanofluids, respectively, as a function of the volume fraction of nanoparticles. The calculated results are represented by the opened circles and the experimental results are also shown with the closed squares for comparison. It is noticed that the experimental results show the anomalous enhancement of thermal conductivity compared with the calculated results. The conventional models, such as H–C model cannot predict thermal conductivity of nanofluids, which means that thermal transport in nanofluids cannot be understood in view of the effective medium theory. New comprehensive theories are needed for the prediction of thermal conductivity of nanofluids. Since it has been reported experimentally that thermal conductivity of nanofluids is related with several factors, such as the stability of suspension of nanoparticles, nanoparticle size, and viscosity of basefluids, the new theories should include the effects of those factors. We presume that thermal conductivity of nanofluids is determined by the compensation of the effects of those factors. More experimental study is needed to provide insight into the mechanism of thermal transport in nanofluids.

4. Conclusion

Thermal conductivities of TiO_2 , Al_2O_3 , Fe, and WO_3 nanofluids were investigated in this study. We found the followings from the comparison of thermal conductivities of the nanofluids: the surface-to-volume ratio of nanoparticles is a principle factor in determining thermal conductivity of nanofluids. Even though the effects of a large number of factors are compensated for thermal conductivity of nanofluids, the intrinsic thermal conductivity of the suspended nanoparticles does not give a primary effect in determining thermal conductivity of nanofluids.

The anomalous enhancement of thermal conductivity of nanofluids could not be understood with the H–C model. New comprehensive theories are needed for the prediction of thermal conductivity of nanofluids. The new theories should include the effects of the factors, such as the stability of suspension of nanoparticles, nanoparticle size, and viscosity of basefluids.

Fundamental understanding of the thermal transport behavior of nanofluids will develop a new generation of highly efficient heat-transfer fluids. Large improvement of the thermal conductivity of nanofluids holds a significant potential for revolutionizing industries that are dependent on the performance of heat transfer fluids. More experimental study should be performed to provide insight into the mechanism of thermal transport in nanofluids.

Acknowledgements

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