# Is It Important to Measure the Volumetric Mass Density of Nanofluids? 

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#### Abstract

The present study aims to measure the volumetric mass density of NiPd-heptane nanofluids synthesized using a one step method known as thermal decomposition of metal-surfactant complexes. The particle concentration is up to $7.55 \mathrm{~g} / 1$ and the temperature range of the experiment is from $20^{\circ} \mathrm{C}$ to $50^{\circ} \mathrm{C}$. The measured values were compared with the mixture theory and good agreement between the theoretical equation and measurement were obtained. Moreover, the available nanofluids volumetric mass density data in the literature is reviewed.


Keywords-NiPd nanoparticles, nanofluids, volumetric mass density, stability.

## I. Introduction

PREPARATION of stable nanofluids is an unavoidable issue that can alter the thermo-physical properties of nanofluids. The stability of nanofluids refers to several aspects: 1) Nanofluids are multi-phase dispersion system with high surface energies and are, therefore, thermodynamic unstable. 2) Nanoparticles dispersed in the nanofluids have strong Brownian motions. The mobility of the nanoparticles can offset their sedimentation caused by the gravity field. 3) Dispersion of nanoparticles in the fluids may deteriorate with time due to the aggregation of nanoparticles, which is caused by van der Waals forces. 4) No chemical reactions either between the suspended nanoparticles or between the base fluid and nanoparticles are desired at working conditions of the nanofluid. Therefore, there are two phenomena that are critical to the stability of nanofluid, aggregation and sedimentation.

In recent years, considerable efforts have been made to prepare stable and durable nanofluids, before it can be fully implemented in the industrial applications [1]-[4]. The nanoparticles that have been mostly used to prepare nanofluids include metallic particles ( $\mathrm{Al}, \mathrm{Au}, \mathrm{Ag}$ and Cu ) and nonmetallic particles $\left(\mathrm{Al}_{2} \mathrm{O}_{3}, \mathrm{ZnO}, \mathrm{Cuo}, \mathrm{TiO}_{2}, \mathrm{Fe}_{3} \mathrm{O}_{4}, \mathrm{CNTs}, \mathrm{SiO}_{2}\right.$ and AIN), with an average size varied between 10 and 80 nm . The base fluids commonly used are water and ethylene glycol.

Although research works have shown that changing the pH of suspensions or adding surfactants can prevent coalescence and sedimentation of nanoparticle, the obtained stability was

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only for few months [4]. The best way to produce long-term stable nanofluids that exhibits true 'nano' behavior, and measure the thermo-physical properties accurately may be a single-step method. In this way, the chance of agglomeration is reduced. Also, reducing the particle size should remarkably decrease the sedimentation speed of the nanoparticles and improve the stability of nanofluids, since the particle's sedimentation velocity is proportional to the square of the particle size. In this context, a bimetallic nanofluid with high stability is produced.

The purpose of this work is twofold. First, synthesize nickel/palladium (NiPd) nanoparticles by thermal decomposition of metal-surfactant complexes. Then, measure the volumetric mass density of NiPd-heptane nanofluids under various operating temperature and nanofluid concentrations, and compared the results with theoretical prediction of mixture theory.

## II. Material and Methods

## A. Nickel-Palladium (Ni-Pd)

Bimetallic PdNi nanoparticles are currently attracting a great deal of interest due to their physical and chemical properties. In particular, it is well-known that Pd is a metal with important existing and potential applications as a catalyst in heterogeneous catalysis. The use of Pd as a single active metal component in catalysis has received considerable attention on the basis of its remarkable activity for oxidation reactions and the availability of cleaner fuels, but at the same time it has negative economical aspects due to its high cost [5]. An interesting approach in these respects consists of alloying Pd with lower cost and higher surface energy metals, since it would be economically attractive to design bimetallic catalyst nanoparticles in which the precious and catalytic Pd atoms segregate to the surface. Promising results in this sense have been obtained by using Ni [5]-[7].

## B. Synthesis of Nanofluids

The synthesis is slightly modified from ref [8]. In a typical experiment, a solution mixture of $1: 1$ nickel acetylacetonate ( Ni (acac)2-0.082g, 0.328 mmol ) and palladium acetylacetonate $\left(\mathrm{Pd}(\mathrm{acac})_{2}-0.1 \mathrm{~g}, 0.328 \mathrm{mmol}\right)$ in 2 mL of trioctylphosphine, was injected into a 3-neck flask containing 7 mL of oleylamine at $205^{\circ} \mathrm{C}$. After aging at $205^{\circ} \mathrm{C}$ for 30 minutes, the temperature was increased to $235^{\circ} \mathrm{C}$ in 30 minutes and kept at this temperature for one hour. Washing of the nanoparticles was done through three cycles using ethanol for flocculation and centrifugation to remove the supernatant. Finally, the nanoparticles were redispersed in pure heptane.

## C. Characterization of Nanoparticles

The morphological characterization of NiPd nanoparticles deposited on a carbon-coated copper grid was carried using a high resolution Transmission Electron Microscopy (HRTEM JEOL 3010).

## D.Formulation of Nanofluids

NiPd-heptane nanofluids with woncentrations of $16 \mathrm{~g} / 1$ and $4 \mathrm{~g} / 1$ were first produced (Fig. 1). Then, five concentrations of $7.55 \mathrm{~g} / 1,5.95 \mathrm{~g} / 1,4.5 \mathrm{~g} / 1,2.95 \mathrm{~g} / 1$ and $1 \mathrm{~g} / \mathrm{l}$ were obtained by dilution with heptane under magnetic stirring for 1 min at 220 rpm (Fisher Scientific/FB15038). The produced samples of 16 $\mathrm{g} / 1$ and $4 \mathrm{~g} / 1$ were stable for one year and a half without any visible sedimentation. It should be noted that the other dilutions were prepared after that period.

Fig. 1 Photograph of NiPd-Heptane Nanofluids

## III. Results and Discussion

## A. Morphology and Crystalline Structure of Nanoparticles

Figs. 2 (a) and (b) show the transmission electron microscopy and high resolution transmission electron microscopy of NiPd nanoparticles, respectively. A high resolution TEM (HRTEM) image of the nanoparticles revealed the highly crystalline nature of the nanoparticles with a diameter of about 5 nm . Moreover, the image shows that nanoparticles are uniform and monodispersed. In Fig. 3, the histogram of TEM image confirms that NiPd nanoparticles have an average diameter of $4.5 \mathrm{~nm}+/-0.6 \mathrm{~nm}$ (standard deviation).


Fig. 2 (a) TEM image, (b) HRTEM image


Fig. 3 Histogram of TEM image of NiPd nanoparticles.

## B. Volumetric Mass Density Measurement

Nanofluid volumetric mass density is measured using the portable DMA 35 according to the oscillating U-tube principle. The sample is introduced into a U-shaped borosilicate glass tube that is being excited to vibrate at its characteristic frequency electronically. The characteristic frequency changes depending on the volumetric mass density of the sample. Through determination of the characteristic frequency the sample can be calculated. Due to the temperature dependency of the density value, the temperature of the sample has to be determined precisely.

To validate the measurements, the volumetric mass density of distilled water was measured at four temperatures $20^{\circ} \mathrm{C}$, $30^{\circ} \mathrm{C}, 40^{\circ} \mathrm{C}$, and $50^{\circ} \mathrm{C}$. The results were compared with the data provided by A. Bejan [9]. Table I shows the measured and reference values. It can be seen that the deviation between measured data and standard reference for the limited temperature range considered in this study is less than $0.2 \%$.

TABLE I

| VOLUMETRIC MASS DENSITY OF DISTILLED WATER |  |  |  |
| :---: | :---: | :---: | :---: |
| $\mathrm{T}\left({ }^{\circ} \mathrm{C}\right)$ | $\rho_{\text {meas }}\left(\mathrm{g} / \mathrm{cm}^{3}\right)$ | $\rho_{\text {ref }}\left(\mathrm{g} / \mathrm{cm}^{3}\right)$ | $\left(\rho_{\text {meas }}-\rho_{\text {reff }} / \rho_{\text {ref }}(\%)\right.$ |
| 20 | 0.9992 | 0.9982 | 0.10 |
| 30 | 0.9967 | 0.9952 | 0.15 |
| 40 | 0.9925 | 0.9922 | 0.03 |
| 50 | 0.9882 | 0.9877 | 0.051 |

## C. Volumetric Mass Density of Nanofluids

Because density is an important property of nanofluids that affects the Reynolds number, friction factors, pressure loss and Nusselt number, we focus our attention in this paper on accurately measuring the density of NiPd-heptane nanofluids. The equation for the density of two-phase mixtures for particles of micrometer size is available in the literature on slurry flows [10]. Pak and Cho [11] measured the volumetric mass density of $\gamma-\mathrm{Al}_{2} \mathrm{O}_{3}$ and $\mathrm{TiO}_{2}$ at a volume concentration of $4.5 \%$. They confirmed that the mixture theory of micrometer particle is valid for nanometer size particles, which is expressed by the following expression

$$
\begin{equation*}
\rho_{n f}=(1-\phi) \rho_{f}+\phi \rho_{s} \tag{1}
\end{equation*}
$$

where $\rho_{n f}$ is the density of the nanofluid, $\rho_{s}$ is the density of the particle, $\phi$ is the particle volume concentration, $\rho_{f}$ and is the density of the base fluid.

Fig. 4 displays the measured volumetric mass density of NiPd-heptane nanofluids for different concentrations and within a temperature range of $20^{\circ} \mathrm{C}$ to $50^{\circ} \mathrm{C}$. The general observation is that volumetric density decreases with temperature and increases with an increase of particle concentrations, which agree well with the behavior of pure and dispersed fluids. Comparison of the measured volumetric mass density values and those calculated from (1) are also presented. The maximum deviation is $1.12 \%$ at a particle concentration of $5.95 \mathrm{~g} / \mathrm{l}$. Therefore, we conclude that mixture theory predicts well the volumetric mass density of nanofluids.


Fig. 4 Volumetric Mass Density of NiPd Nanofluids

Table II summarizes experimental data of nanofluids volumetric mass density reported by previous authors and their deviation from the proposed theoretical model [12]-[19]. Most of experimental findings show that mixture theory predicts well the volumetric mass density of nanofluids, which is not a new finding since this theory is based on the mass conservation. However, only a few authors that reported contradictory results; Kumaresan and Velraj [16] showed that the mixture theory could not predict the volumetric mass density of the nanofluids with the presence of MWCNT nanoparticles. They attributed this difference to the spontaneous filling of nanotubes with water in a confined way which increases the mass of the nanofluid for a given volume. Mahbubul et al. [19] explained this difference to be mainly due to the fact that pack and cho model was derived for water based nanofluids.

## IV. CONCLUSION

Well dispersed heptanes based NiPd nanofluids with average particle size of 4.5 nm were synthesized using thermal decomposition of metal-surfactant complexes. The samples were stable at least one year and a half and there was no visible sedimentation. As results show, the nanofluid volumetric mass density can be determined using mixture theory since it is based on mass conservation. However, the contradictory results reported in literature can be due to the fact that nanofluids can not be prepared precisely at a desired volume or mass concentrations.

TABLE II
Summary of Literature Review of Volumetric Mass
Density of Nanofluids

| DENSITY OF NANOFLUIDS |  |  |
| :---: | :---: | :---: |
| Authors | Nanofluids /parameters | Comparison with theory |
| Ho et al. [12] | $\mathrm{Al}_{2} \mathrm{O}_{3}$-water, $\Phi=0.1-4$ vol. $\%, \mathrm{D}=33 \mathrm{~nm}, \mathrm{~T}=5-40^{\circ} \mathrm{C}$ | Agree well |
| Ravikanth et al. [13] | $\mathrm{Al}_{2} \mathrm{O}_{3}$-water/EG, $\Phi=2-10$ vol. $\%, \mathrm{D}=53 \mathrm{~nm}, \mathrm{~T}=0-50{ }^{\circ} \mathrm{C}$ | Agree well |
| Mohamed Ali et al. [14] | $\mathrm{Al}_{2} \mathrm{O}_{3}$-water, $\Phi=0.21-0.75$ vol. $\%, \mathrm{D}=10 \mathrm{~nm}, \mathrm{~T}=25.5{ }^{\circ} \mathrm{C}$ | Agree well |
| Fakoor Pakdaman et al. [15] | MWCNT-oil, $\Phi=0.1,0.2$ and $0.4 \mathrm{wt} . \%, \mathrm{D} \times \mathrm{H}=2-6 \mathrm{~nm} \times 1-10 \mu \mathrm{~m}$ | satisfactorily agreement |
| Kumaresan and Velraj [16] | CNT-water/EG, $\Phi=0.15,0.3$ and $0.45 \mathrm{wt} \%,. \mathrm{D} \times \mathrm{H}=30-50 \mathrm{~nm} \times 10-20 \mu \mathrm{~m}$ | Overestimated |
| Heyhat et al. [17] | $\mathrm{Al}_{2} \mathrm{O}_{3}$-water, $\Phi=0.1-2 \mathrm{vol} . \%, \mathrm{D}=40 \mathrm{~nm}, \mathrm{~T}=20-60^{\circ} \mathrm{C}$ | Agree well |
| Saeedinia et al. [18] | CuO-oil, $\Phi=0.2-2 \mathrm{wt} . \%, \mathrm{D}=50 \mathrm{~nm}, \mathrm{~T}=20-60^{\circ} \mathrm{C}$ | Agree well |
| Mahbubul et al. [19] | $\mathrm{Al}_{2} \mathrm{O}_{3}-\mathrm{R} 141 \mathrm{~b}, \Phi=0.1-0.4$ vol. $\%, \mathrm{D}=13 \mathrm{~nm}, \mathrm{~T}=5-20^{\circ} \mathrm{C}$ | underestimated |

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