This article was downloaded by: On: *28 January 2011* Access details: *Access Details: Free Access* Publisher *Taylor & Francis* Informa Ltd Registered in England and Wales Registered Number: 1072954 Registered office: Mortimer House, 37-41 Mortimer Street, London W1T 3JH, UK



## Physics and Chemistry of Liquids

Publication details, including instructions for authors and subscription information: http://www.informaworld.com/smpp/title~content=t713646857

# Density and Viscosity of Methanol Trichloroethylene, *n*-Propanol Trichloroethylene and *n*-Butanol Trichloroethylene Mixtures

**To cite this Article** (2000) 'Density and Viscosity of Methanol Trichloroethylene, *n*-Propanol Trichloroethylene and *n*-Butanol Trichloroethylene Mixtures', Physics and Chemistry of Liquids, 38: 4, 433 - 438

**To link to this Article: DOI:** 10.1080/00319100008030290

**URL:** http://dx.doi.org/10.1080/00319100008030290

# PLEASE SCROLL DOWN FOR ARTICLE

Full terms and conditions of use: http://www.informaworld.com/terms-and-conditions-of-access.pdf

This article may be used for research, teaching and private study purposes. Any substantial or systematic reproduction, re-distribution, re-selling, loan or sub-licensing, systematic supply or distribution in any form to anyone is expressly forbidden.

The publisher does not give any warranty express or implied or make any representation that the contents will be complete or accurate or up to date. The accuracy of any instructions, formulae and drug doese should be independently verified with primary sources. The publisher shall not be liable for any loss, actions, claims, proceedings, demand or costs or damages whatsoever or howsoever caused arising directly or indirectly in connection with or arising out of the use of this material.

*Phys. Chem. Liq.*, 2000, Vol. 38, pp. 433-438 Reprints available directly from the publisher Photocopying permitted by license only

# DENSITY AND VISCOSITY OF METHANOL + TRICHLOROETHYLENE, *n*-PROPANOL + TRICHLOROETHYLENE AND *n*-BUTANOL + TRICHLOROETHYLENE MIXTURES

### T. E. VITTAL PRASAD<sup>a</sup>, S. B. AGRAWAL<sup>a, b</sup>, A. B. BAJAJ<sup>a, b</sup> and D. H. L. PRASAD<sup>a</sup>, \*

<sup>a</sup> Properties Group, Chemical Engineering Laboratory, Indian Institute of Chemical Technology, Hyderabad-500 007, India; <sup>b</sup> College of Engineering and Technology, Akola, India

(Received 4 May 1999)

Density and Viscosity measurements on the binary mixtures of methanol + trichloroethylene, n-propanol + trichloroethylene, and n-butanol + trichloroethylene binary mixtures at 303.15, 313.15 and 323.15K are reported. The representation of the data by simple mixing rules is also studied.

Keywords: Density; viscosity; mixtures; methanol; n-propanol; n-butanol; trichloroethylene

#### INTRODUCTION

In continuation of our investigations on the thermophysical properties of the binary systems formed by an aliphatic alcohol (methanol, ethanol, *n*-propanol, *n*-butanol) as one component and chloroethane (1,2-dichloroethane, 1,1,1-trichloroethane, 1,1,2,2-tetrachloroethane) as the other component [1-6], this study on the density and viscosity of the binary mixtures noted in the title has been carried out. No published data on these mixtures is available for comparison.

<sup>\*</sup>Corresponding author.

#### **EXPERIMENTAL**

#### Materials

Spectrographic grade methanol procured from SDS Fine Chemicals, Boisar, India has been distilled twice after drying over silica gel. The middle fraction of the second distillation (boiling at 336.45 K, corresponding to the local atmospheric pressure) has been stored in amber coloured bottles for use in the experiments.

n-Propanol supplied by E. Merck, Bombay, India is purified by adding a small quantity of sodium metal and refluxing for 2 h. followed by treatment with anhydrous sodium carbonate, drying over calcium hydroxide and distillation.

A R grade *n*-butanol procured from SDS Fine Chemicals, Boisar, India has been purified by washing with dilute sulfuric acid and sodium bisulfite solution to remove the bases, aldehydes, ketones. Esters are removed by boiling for 1.5 h with 20% sodium hydroxide solution. The alcohol is dried with potassium carbonate followed by barium oxide. The final purification is by means of careful distillation.

Trichloroethylene supplied by Ranbaxy Laboratories, SAS Nagar, Punjab, India is treated with potassium carbonate solution followed by water. After drying over potassium carbonate and calcium chloride trichloroethylene distilled twice, each time collecting the middle 40% fraction.

#### Methods

Measurements on the density of the pure liquids as well as the mixtures are carried out using pycnometers, carefully calibrated by weighing double distilled water. The weights required in the entire work are recorded using a Mettler balance, accurate to  $\pm 0.0001$  g. The temperatures are maintained to within  $\pm 0.05$  K, of the desired value by keeping the pycnometers in an electronically controlled water bath for sufficient length of time (usually 1h).

A Haake falling ball viscometer with provision to maintain the temperature of the test liquid to within  $\pm 0.05$  K of the desired value, is used to determine the viscosity of the pure liquids and mixtures. When the experiments are conducted, care is taken to avoid entrainment of any air bubbles or particulate matter in the apparatus. Care is also

Substance	Refract	tive-index	Density, g/mL		
	This work	Literature [8]	This work	Literature [7]	
Methanol	1.3288	1.3288	0.7911	0.7910	
n-Propanol	1.3489	1.3850	0.8041	0.8040	
n-Butanol	1.3994	1.3993	0.8101	0.8100	
Trichloroethylene	1.4764	1.4765	1.4621	1.4620	

TABLE I Comparison of Refractive-index and Density of the substances used in the present work with Literature data at  $293.15 \, \text{K}$ 

taken to avoid parallax error. The time measurements are accurate to  $\pm 0.05$  s. Based on the comparison of the data on pure liquids with the literature data presented in Table I, and several other measurements, collected earlier with the literature data, the measurements of the viscosity reported in this paper are expected to be within  $\pm 0.5\%$ .

Mixture samples are prepared from weighed quantities of the pure liquids and the constancy of the composition is checked after each experiment by gas chromatographic analysis. All the measurements have been carried out three times. The triplicate measurements are within  $\pm 0.2\%$  of the values given in Tables III-V.

#### **RESULTS AND DISCUSSION**

The densities and viscosities measured in the present work are presented in Tables II-V. The law of additive volumes, rearranged for convinence as

$$\rho_m = \rho_1 \rho_2 / (\rho_2 x_1 + \rho_1 x_2) \tag{1}$$

represents all the mixture density measurements of the present work with an average absolute deviation of 0.5%. Linear law

$$\eta_m = \eta_1 x_1 + \eta_2 x_2 \tag{2}$$

Arrhenius equation

$$\eta_m = \mathbf{e}^{(x_1 \ln \eta_1 + x_2 \ln \eta_2)} \tag{3}$$

Substance	Viscosity, cP			
	This work	Literature		
Methanol	0.5944	0.5945		
n-Propanol	2.1925	2.2000		
n-Butanol	2.8749	2.8750		
Trichloroethylene	0.6219	0.6220		

TABLE II Comparison of the pure liquid viscosity data at 293.15 K with Literature data from Ref. [9]

TABLE III Density and viscosity of methanol + trichloroethylene mixtures

Mole fraction of methanol	Temperature Density, g/mL	e = 303.15 K Viscosity, cP	Temperature Density, g/mL	e = 313.15 K Viscosity, cP	Temperatur Density, g/mL	e = 323.15 K Viscosity, cP
1.0000	0.7910	0.5020	0.7760	0.4870	0.7660	0.4770
0.9011	0.8560	0.5080	0.8400	0.4920	0.8290	0.4760
0.7560	0.9520	0.5180	0.9360	0.4970	0.9230	0.4750
0.5963	1.0570	0.5300	1.0400	0.5050	1.0260	0.4740
0.4460	1.1562	0.5400	1.1382	0.5110	1.2320	0.4730
0.3071	1.2480	0.5500	1.2294	0.5180	1.2130	0.4735
0.1523	1.3510	0.5600	1.3390	0.5240	1.3240	0.4730
0.000	1.4620	0.5710	1.4420	0.5310	1.4230	0.4720

TABLE IV Density and viscosity of n-propanol + trichloroethylene mixtures

Mole fraction of n-Propanol	Temperatur Density, g/mL	e = 303.15 K Viscosity, cP	Temperatur Density, g/mL	e = 313.15 K Viscosity, cP	Temperatur Density, g/mL	e = 323.15 K Viscosity, cP
1.0000	0.7920	1.7020	0.7820	1.3360	0.7720	1.1110
0.9543	0.8890	1.5370	0.8880	1.2190	0.8780	1.0180
0.6998	0.9930	1.3700	0.9800	1.0950	0.9680	0.9190
0.5454	1.0960	1.1930	1.0810	0.9700	1.0680	0.8200
0.3927	1.1980	1.0190	1.1820	0.8480	1.1670	0.7220
0.2530	1.2910	0.8600	1.2740	0.7350	1.2580	0.6340
0.1062	1.3910	0.6920	1.3720	0.6170	1.3530	0.5390
0.0240	1.4450	0.5980	1.4260	0.5500	1.4050	0.4870
0.0000	1.4620	0.5710	1.4420	0.5310	1.4230	0.4720

TABLE V Density and viscosity of n-butanol + trichloroethylene mixtures

Mole fraction on n-Butanol	Temperatur Density, g/mL	e = 303.15 K Viscosity, cP	Temperatur Density, g/mL	e = 313.15 K Viscosity, cP	Temperatur Density, g/mL	e = 323.15 K Viscosity, cP
1.0000	0.7980	2.2620	0.7875	1.7920	0.7825	1.3290
0.8476	0.8980	2.0005	0.8870	1.5990	0.8790	1.2520
0.6960	0.9993	1.7470	0.9860	1.4080	0.9760	1.1110
0.5452	1.0990	1.4920	1.0840	1.2180	1.0730	0.9730
0.3954	1.1990	1.2400	1.1810	1.0300	1.1690	0.8350
0.2464	1.3000	0.9800	1.2800	0.8140	1.2640	0.6980
0.0983	1.3720	0.6690	1.3530	0.6010	1.3350	0.5200
0.0000	1.4620	0.5710	1.4420	0.5310	1.4230	0.4720

System	Number of		Percent average absolute deviation				
	data points	Linear law	Arrhenius equation	Kendall and Monroe equation	Lobe's equation		
Methanol +							
Trichloroethylene <i>n</i> -Propanol +	24	0.47	0.57	0.90	1.98		
Trichloroethylene $n$ -Butanol +	27	0.22	0.24	1.05	1.81		
Trichloroethylene	24	0.14	0.15	1.31	1.65		
Overall	75	0.28	0.32	1.09	1.81		

TABLE VI Representation of the viscosity data by simple mixing rules

Kendall and Monroe equation

$$\eta_m = \left(x_1 \eta_1^{1/3} + x_2 \eta_2^{1/3}\right)^3 \tag{4}$$

and Lobe's equation

$$(\eta_m/\rho_m) - \Phi_1(\eta_1/\rho_1) e^{\Phi_2\beta} + \Phi_2(\eta_2/\rho_2) e^{\Phi_1\alpha}$$
(5)

with

$$\alpha = -1.7 \ln(\eta_2/\rho_2)/(\eta_1/\rho_1)$$
(6)

$$\beta = 0.27 \ln \left[ \frac{(\eta_2/\rho_2)}{(\eta_1/\rho_1)} \right] + 1.3 \ln \left[ \frac{(n_2/\rho_2)}{(n_1/\rho_1)} \right]^{0.5}$$
(7)

have been used to calculate the mixture viscosity. A summary of the comparisons given in Table VI shows that on an overall basis the linear law is preferable with an average absolute deviation of 0.28%, comparable to the Arrhenius equation with 0.32%, Kendall and Monroe equation with 1.09% and Lobe's equation with 1.81%.

#### References

- Srinivas Rao, G., Malla Reddy, V. and Prasad, D. H. L. (1989). Physics and Chemistry of Liquids, 20, 87.
- [2] Sivaram Prasad, G., Venkateshwara Rao, M. and Prasad, D. H. L. (1990). Journal of Chemical and Engineering Data, 35, 122.
- [3] Srinivas, Ch., Venkateshwara Rao, M. and Prasad, D. H. L. (1991). Fluid Phase Equilibria, 69, 285
- [4] Lakshman, V., Venkateshwara Rao, M. and Prasad, D. H. L. (1991). Fluid Phase Equilibria, 69, 270.

- [5] Kiran Kumar, R., Venkateshwara Rao, M. and Prasad, D. H. L. (1991). Fluid Phase Equilibria, 70, 19.
- [6] Jaya Prakash, D., Sree Lakshmi, D., Venkateshwara Rao, M. and Prasad, D. H. L. (1996). Physics and Chemistry of Liquids, 33, 249.
- [7] Reid, R. C., Prausnitz, J. M. and Sherwood, T. K. (1977). The Properties of Gases and Liquids, Third Edition, Mc GrawHill, New York.
- [8] Weast, R. C. and Astle, M. J. (1985). CRC Handbook of Data on Organic Compounds, CRC Press, Boca Raton, Florida.
- [9] Gallant, R. W., Physical Properties of Hydrocarbons, Vols. I & II, Gulf Publishing Co., Houston, Texas.