

COMPONENTS: (1) 1-Heptene; C ₇ H ₁₄ ; [592-76-7] (2) Water; H ₂ O; [7732-18-5]	EVALUATOR: G.T. Hefter, School of Mathematical and Physical Sciences, Murdoch University, Perth, W.A., Australia. July 1985
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CRITICAL EVALUATION:

Quantitative data for the solubility of water (2) in 1-heptene (1) have been reported in the publications listed in Table 1.

TABLE 1: Quantative Solubility Studies of
Water (2) in 1-Heptene (1)

Reference	T/K	Method
Black <i>et al.</i> (ref 1)	283-294	radiotracer
Englin <i>et al.</i> (ref 2)	283-303	analytical

The solubility of 1-heptene in dilute aqueous HNO₃ solutions have also been reported (ref 3). However, as no other data are available for comparison these values will not be considered in this Evaluation. The interested user is referred to the relevant Data Sheet immediately following this Critical Evaluation.

The reported values (ref 1 and 2) for the solubility of water in 1-heptene are listed in Table 2.

The data are in very poor agreement. In the absence of confirmatory studies selection of best values is difficult. However, it may be noted that in well characterized systems (e.g., water in benzene) the results of Englin *et al.* at $T < 300\text{K}$ are generally reliable whilst those of Black *et al.* generally differ markedly from Recommended values. Thus, values of Englin *et al.* (ref 2) have been selected as the "Best" values available although, in the absence of confirmatory studies, it is not possible at this stage to reject outright the data of Black *et al.* (ref 1).

(continued next page)

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CRITICAL EVALUATION: (continued)

TABLE 2: Tentative Values of the Solubility
of Water (2) in 1-Heptene (1)

T/K	Solubility Values		
	Reported values 10 ² g(2)/100g sln	"Best" values ^a 10 ² g(2)/100g sln	10 ³ x ₂
283	6.92 (ref 1), 1.86 (ref 2)	1.9	1.0
293	11.2 (ref 1), 2.49 (ref 2)	2.5	1.4
298	3.2 ^b (ref 2)	3.2	1.7
303	3.75 (ref 2)	3.8	2.0

^aRounded values of Englin *et al.* (ref 2); see text.

^bObtained by the Evaluator by graphical interpolation.

REFERENCES

- Black, C.; Joris, G.G.; Taylor, H.S. *J. Chem. Phys.* 1948, *16*, 537-43.
- Englin, B.A.; Plate, A.F.; Tugolukov, V.M.; Pryanishnikova, M.A. *Khim. Tekhnol. Topl. Masel* 1965, *10*, 42-6.
- Natarajan, G.S.; Venkatachalam, K.A. *J. Chem. Eng. Data.* 1972, *17*, 328-9.

COMPONENTS: (1) 1-Heptene; C_7H_{14} ; [592-76-7] (2) Water; H_2O ; [7732-18-5]	ORIGINAL MEASUREMENTS: Black, C.; Joris, G.G.; Taylor, H.S. <i>J. Chem. Phys.</i> <u>1948</u> , <i>16</i> , 537-43.																				
VARIABLES: Temperature: 10 - 21.2°C	PREPARED BY: A. Maczynski and Z. Maczynska																				
EXPERIMENTAL VALUES: <p style="text-align: center;">Solubility of water in 1-heptene</p> <table border="1" style="width: 100%; border-collapse: collapse;"> <thead> <tr> <th style="text-align: center;">$t/^\circ C$</th> <th style="text-align: center;">$g(2)/100\ g(1)$</th> <th style="text-align: center;">$g(2)/100\ g\ sln$ (compiler)</th> <th style="text-align: center;">$10^3 x_2$ (compiler)</th> </tr> </thead> <tbody> <tr> <td style="text-align: center;">10.0</td> <td style="text-align: center;">0.0692</td> <td style="text-align: center;">0.0692</td> <td style="text-align: center;">3.76</td> </tr> <tr> <td style="text-align: center;">20.1</td> <td style="text-align: center;">0.1126</td> <td style="text-align: center;">0.1126</td> <td style="text-align: center;">6.11</td> </tr> <tr> <td style="text-align: center;">20.5</td> <td style="text-align: center;">0.1047</td> <td style="text-align: center;">0.1047</td> <td style="text-align: center;">5.68</td> </tr> <tr> <td style="text-align: center;">21.2</td> <td style="text-align: center;">0.1158</td> <td style="text-align: center;">0.1158</td> <td style="text-align: center;">6.28</td> </tr> </tbody> </table> <p style="text-align: center;">(at total saturation pressure of 1 atm)</p>		$t/^\circ C$	$g(2)/100\ g(1)$	$g(2)/100\ g\ sln$ (compiler)	$10^3 x_2$ (compiler)	10.0	0.0692	0.0692	3.76	20.1	0.1126	0.1126	6.11	20.5	0.1047	0.1047	5.68	21.2	0.1158	0.1158	6.28
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AUXILIARY INFORMATION																					
METHOD/APPARATUS/PROCEDURE: Air saturated with radioactive water vapor was bubbled through (1) until saturation was attained. Dissolved water was separated from (1) by absorption on calcium oxide. The tritium was transferred into the counter by equilibration with ethanol vapor. The method is described in ref (1).	SOURCE AND PURITY OF MATERIALS: (1) Havard University; purity not specified; used as received. (2) not specified. ESTIMATED ERROR: soly. \pm 1% (type or error not specified) REFERENCES: 1. Joris, G.G.; Taylor, H.S. <i>J. Chem. Phys.</i> <u>1948</u> , <i>16</i> , 45.																				

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VARIABLES: Temperature: 10-30°C	PREPARED BY: A. Maczynski and Z. Maczynska												
EXPERIMENTAL VALUES: <p style="text-align: center;">Solubility of water in 1-heptene</p> <table border="1" style="margin-left: auto; margin-right: auto;"> <thead> <tr> <th style="text-align: center;"><u>t/°C</u></th> <th style="text-align: center;"><u>g(2)/100 g sln</u></th> <th style="text-align: center;"><u>10³x₂ (compiler)</u></th> </tr> </thead> <tbody> <tr> <td style="text-align: center;">10</td> <td style="text-align: center;">0.0186</td> <td style="text-align: center;">1.01</td> </tr> <tr> <td style="text-align: center;">20</td> <td style="text-align: center;">0.0249</td> <td style="text-align: center;">1.36</td> </tr> <tr> <td style="text-align: center;">30</td> <td style="text-align: center;">0.0375</td> <td style="text-align: center;">2.04</td> </tr> </tbody> </table>		<u>t/°C</u>	<u>g(2)/100 g sln</u>	<u>10³x₂ (compiler)</u>	10	0.0186	1.01	20	0.0249	1.36	30	0.0375	2.04
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10	0.0186	1.01											
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METHOD/APPARATUS/PROCEDURE: Component (1) was introduced into a thermostatted flask and saturated for 5 hr. with (2). Next, calcium hydride was added and the evolving hydrogen volume measured and hence the concentration of (2) in (1) was evaluated.	SOURCE AND PURITY OF MATERIALS: (1) not specified. (2) not specified. ESTIMATED ERROR: Not specified. REFERENCES:												

COMPONENTS: (1) 1-Heptene; C ₇ H ₁₄ ; [592-76-7] (2) Water; H ₂ O; [7732-18-5]	ORIGINAL MEASUREMENTS: Natarajan, G.S.; Venkatachalam, K.A. <i>J. Chem. Eng. Data</i> <u>1972</u> , 17, 328-9																
VARIABLES: Temperature: 20-30°C	PREPARED BY: M.C. Haulait-Pirson, G.T. Hefter																
EXPERIMENTAL VALUES: Solubility of 1-heptene in 0.001 mol L ⁻¹ HNO ₃ solution <table border="1" data-bbox="207 516 1126 741"> <thead> <tr> <th>t/°C</th> <th>10⁴ mol L⁻¹ sln^a</th> <th>10³ g(1)/100 g sln^b (compiler)</th> <th>10⁶ x₁ (compiler)</th> </tr> </thead> <tbody> <tr> <td>20</td> <td>3.16 ± 0.13</td> <td>3.1</td> <td>5.7</td> </tr> <tr> <td>25</td> <td>2.81 ± 0.12</td> <td>2.8</td> <td>5.1</td> </tr> <tr> <td>30</td> <td>2.45 ± 0.15</td> <td>2.4</td> <td>4.4</td> </tr> </tbody> </table> <p data-bbox="134 862 1081 923"> a Uncertainties stated to be "standard deviations from means". b Assuming a solution density of 1.00 g mL⁻¹ at all temperatures. </p> <p data-bbox="134 983 1215 1124"> <u>Compiler's note:</u> Although the data have not been measured in pure water the low concentration of the added acid is unlikely to cause the olefin solubility to differ markedly from that in pure water. Further solubility data are given in the paper for 0.05 and 0.1 mol L⁻¹ HCl . </p>		t/°C	10 ⁴ mol L ⁻¹ sln ^a	10 ³ g(1)/100 g sln ^b (compiler)	10 ⁶ x ₁ (compiler)	20	3.16 ± 0.13	3.1	5.7	25	2.81 ± 0.12	2.8	5.1	30	2.45 ± 0.15	2.4	4.4
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METHOD/APPARATUS/PROCEDURE: 15 mL of the aqueous medium was equilibrated with 1 mL of (1) by mechanical shaking in a thermostatted glass burette. After settling (judged visually), 5 mL of the aqueous layer was withdrawn and the olefin content determined by titration with bromine using standard procedures.	SOURCE AND PURITY OF MATERIALS: (1) Matheson, Coleman and Bell; 99%. (2) Not specified. <table border="1" data-bbox="677 1588 1222 1721"> <tbody> <tr> <td> ESTIMATED ERROR: Temp. ± 0.05 K Soly. see table above. </td> </tr> <tr> <td> REFERENCES: </td> </tr> </tbody> </table>	ESTIMATED ERROR: Temp. ± 0.05 K Soly. see table above.	REFERENCES:														
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