

<p>COMPONENTS:</p> <p>(1) 1,2-Difluorobenzene; C₆H₄F₂; [367-11-3]</p> <p>(2) Water; H₂O; [7732-18-5]</p>	<p>EVALUATOR:</p> <p>A. L. Horvath, Imperial Chemical Industries Limited, Runcorn, England.</p> <p>May 1979.</p>								
<p>CRITICAL EVALUATION:</p> <p>The experimental results consist of a single value at 298.15 K published in 1979 by Yalkowsky et al. (1) of the Upjohn Company, Kalamazoo, Michigan. According to the authors (2), the accuracy of the experimental determinations was ± 10 percent. This is a reasonable observation considering the details available on the experimental procedure.</p> <p>The manufacturer's reagent (Aldrich and Eastman) was neither further purified nor degassed before use. The time required for equilibration was indicated as between 4 and 48 hours. The saturated solutions were assayed spectrophotometrically.</p> <p>The solubility value was the average of at least two independent determinations according to the investigators. The experimental result was expressed in Briggsian logarithms only with three significant figures.</p> <p>The objective of the solubility measurements was to extend the correlation technique of aqueous solubilities to a broader group of planar nonelectrolytes by involving the melting points and total molecular surface areas as dependent variables. Using such methods, the authors also showed that branched and cyclic compounds have greater solubilities in water than corresponding linear compounds.</p> <p>To compare, relate, and correlate the solubility data for liquid halogenated benzenes in water, use was made of the theoretical relationship between the molar solubility and solute molar volumes at 298.15 K as discussed in the Introduction. A data comparison with previously selected solubilities shows that new measurements are urgently required in order to resolve the anomalous trend in the correlated experimental data reported by Yalkowsky et al.</p> <p>The uncertainty in the single reported solubility value may be as large as 10 percent or even larger. The following solubility value for 1,2-difluorobenzene in water is tentative:</p> <table border="1" data-bbox="266 1058 900 1139"> <thead> <tr> <th>T/K</th> <th>$10^2 \text{ mol}(1)/\text{dm}^3$</th> <th>g(1)/kg</th> <th>$10^4 x(1)$</th> </tr> </thead> <tbody> <tr> <td>298.15</td> <td>1.0</td> <td>1.14</td> <td>1.81</td> </tr> </tbody> </table> <p style="text-align: center;">REFERENCES</p> <ol style="list-style-type: none"> 1. Yalkowsky, S.H.; Orr, R. J.; Valvani, S. C. <i>Ind. Eng. Chem. Fundam.</i> <u>1979</u>, <i>18</i>(4), 351-3. 2. Yalkowsky, S.H., Personal Communication, <u>1979</u>. 		T/K	$10^2 \text{ mol}(1)/\text{dm}^3$	g(1)/kg	$10^4 x(1)$	298.15	1.0	1.14	1.81
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COMPONENTS: (1) 1,2-Difluorobenzene; $C_6H_4F_2$; [367-11-3] (2) Water; H_2O ; [7732-18-5]	ORIGINAL MEASUREMENTS: Yalkowsky, S. H.; Orr, R. J.; Valvani, S. C. <i>Ind. Eng. Chem. Fundam.</i> <u>1979</u> , <i>18</i> (4), 351-3.								
VARIABLES: One temperature	PREPARED BY: A. L. Horvath								
EXPERIMENTAL VALUES: <table data-bbox="184 466 865 558" style="width: 100%; border-collapse: collapse;"> <thead> <tr> <th style="text-align: left;">$t/^\circ C$</th> <th style="text-align: left;">$g(l)/dm^3$ ^a</th> <th style="text-align: left;">$10^2 mol(l)/dm^3$ ^b</th> <th style="text-align: left;">$10^4 x(l)$ ^c</th> </tr> </thead> <tbody> <tr> <td style="text-align: left;">25</td> <td style="text-align: left;">1.141</td> <td style="text-align: left;">1.00</td> <td style="text-align: left;">1.809</td> </tr> </tbody> </table> <p data-bbox="184 603 559 676" style="margin-left: 20px;"> a. Calculated by F. W. Getzen. b. Reported. c. Calculated by compiler. </p>		$t/^\circ C$	$g(l)/dm^3$ ^a	$10^2 mol(l)/dm^3$ ^b	$10^4 x(l)$ ^c	25	1.141	1.00	1.809
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AUXILIARY INFORMATION									
METHOD/APPARATUS/PROCEDURE: A small excess of 1,2-difluorobenzene in water was agitated at room temperature for a period of about 24 hours and then filtered. The filtrate was diluted and assayed spectrophotometrically. The determination was done in duplicate.	SOURCE AND PURITY OF MATERIALS: $C_6H_4F_2$: Aldrich commercial grade, used as received. H_2O : Deionized. <table data-bbox="672 1538 1227 1666" style="width: 100%; border-collapse: collapse; margin-top: 10px;"> <tr> <td colspan="2">ESTIMATED ERROR:</td> </tr> <tr> <td>Solubility:</td> <td>$\pm 10\%$.</td> </tr> <tr> <td>Temperature:</td> <td>± 1 K.</td> </tr> </table> REFERENCES:	ESTIMATED ERROR:		Solubility:	$\pm 10\%$.	Temperature:	± 1 K.		
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