

COMPONENTS: (1) Praseodymium bromide; PrBr_3 ; [13536-53-3] (2) 1,2-Diethoxyethane; $\text{C}_6\text{H}_{14}\text{O}_2$; [629-14-1]	ORIGINAL MEASUREMENTS: Kirmse, E. M. <i>Tr. II Vses. Konf. po Teor. Rastvorov</i> 1971, 200-6.
VARIABLES: T/K = 298	PREPARED BY: T. Mioduski and M. Salomon
EXPERIMENTAL VALUES: <p>The solubility of PrBr_3 in 1,2-diethoxyethane at 25°C was given as</p> <p style="text-align: center;">0.6 mass %</p> <p>The corresponding value of the molality calculated by the compiler is</p> <p style="text-align: center;">0.016 mol kg^{-1}</p> <p>The nature of the solid phase was not specified.</p>	
AUXILIARY INFORMATION	
METHOD/APPARATUS/PROCEDURE: Experimental details not given, but were probably similar to previous works of the author which are compiled throughout this volume.	SOURCE AND PURITY OF MATERIALS: Nothing specified, but based on previous work by the author the anhydrous salt was probably prepared by the method of Taylor and Carter (1). ESTIMATED ERROR: Nothing specified. REFERENCES: 1. Taylor, M.D.; Carter, C.P. <i>J. Inorg. Nucl. Chem.</i> 1962, 24, 387.

<p>COMPONENTS:</p> <p>(1) Praseodymium bromide; PrBr_3; [13536-53-3]</p> <p>(2) 1,4-Dioxane; $\text{C}_4\text{H}_8\text{O}_2$; [123-91-1]</p>	<p>ORIGINAL MEASUREMENTS:</p> <p>Kirmse, E.M.; Zwietasch, K.J.; Tirschmann, J.; Oelsner, L.; Niedergesaess, U. <i>Z. Chem.</i> <u>1968</u>, <i>8</i>, 472-3.</p> <p>Kirmse, E.M. <i>Tr. II Vses. Konf. po Teor. Rastvorov.</i> <u>1971</u>, 200-6.</p>
<p>VARIABLES:</p> <p>Room temperature: T/K around 298</p>	<p>PREPARED BY:</p> <p>T. Mioduski</p>
<p>EXPERIMENTAL VALUES:</p> <p>The solubility of PrBr_3 in p-dioxane at around 25°C was given as</p> <p style="text-align: center;">0.3₅ mass %</p> <p>The corresponding molality calculated by the compiler is</p> <p style="text-align: center;">$9.2 \times 10^{-3} \text{ mol kg}^{-1}$</p> <p>The nature of the solid phase was not specified.</p>	
<p>AUXILIARY INFORMATION</p>	
<p>METHOD/APPARATUS/PROCEDURE:</p> <p>The solute-solvent mixtures were isothermally agitated at 25°C or at room temperature. Authors state that the difference found for the solubility was within experimental error limits.</p> <p>Pr was determined by complexometric titration.</p> <p>No other details given.</p>	<p>SOURCE AND PURITY OF MATERIALS:</p> <p>The anhydrous salt was prepared by the method of Taylor and Carter (1).</p> <p>No other information given.</p> <hr/> <p>ESTIMATED ERROR:</p> <p>Nothing specified.</p> <hr/> <p>REFERENCES:</p> <p>1. Taylor, M.D.; Carter, C.P. <i>J. Inorg. Nucl. Chem.</i> <u>1962</u>, <i>24</i>, 387.</p>

COMPONENTS: (1) Praseodymium bromide; PrBr_3 ; [13536-53-3] (2) Alkyl amines	ORIGINAL MEASUREMENTS: Kirmse, E. M. <i>Tr. II Vses. Kong. po Teor. Rastvorov</i> <u>1971</u> , 200-6.																																			
VARIABLES: T/K = 298	PREPARED BY: T. Mioduski and M. Salomon																																			
EXPERIMENTAL VALUES: <table border="1" style="width: 100%; border-collapse: collapse;"> <thead> <tr> <th colspan="3"></th> <th colspan="2" style="text-align: center;">PrBr_3 solubility^a</th> </tr> <tr> <th style="text-align: left;">solvent</th> <th></th> <th></th> <th style="text-align: center;">mass %</th> <th style="text-align: center;">mol kg⁻¹</th> </tr> </thead> <tbody> <tr> <td>1-propanamine;</td> <td>n-C₃H₉N;</td> <td>[107-10-8]</td> <td style="text-align: center;">21.6</td> <td style="text-align: center;">0.724</td> </tr> <tr> <td>2-propanamine;</td> <td>iso-C₃H₉N;</td> <td>[75-31-0]</td> <td style="text-align: center;">9.6</td> <td style="text-align: center;">0.279</td> </tr> <tr> <td>1-butanamine;</td> <td>n-C₄H₁₁N;</td> <td>[109-73-9]</td> <td style="text-align: center;">9.2</td> <td style="text-align: center;">0.266</td> </tr> <tr> <td>2-butanamine;</td> <td>sec-C₄H₁₁N;</td> <td>[13952-84-6]</td> <td style="text-align: center;">26.5</td> <td style="text-align: center;">0.947</td> </tr> <tr> <td>di-2-butylamine;</td> <td>(sec-C₄H₉)₂NH;</td> <td>[626-23-3]</td> <td style="text-align: center;">0.04</td> <td style="text-align: center;">0.0011</td> </tr> </tbody> </table> <p>^aMolalities calculated by the compilers.</p>					PrBr_3 solubility ^a		solvent			mass %	mol kg ⁻¹	1-propanamine;	n-C ₃ H ₉ N;	[107-10-8]	21.6	0.724	2-propanamine;	iso-C ₃ H ₉ N;	[75-31-0]	9.6	0.279	1-butanamine;	n-C ₄ H ₁₁ N;	[109-73-9]	9.2	0.266	2-butanamine;	sec-C ₄ H ₁₁ N;	[13952-84-6]	26.5	0.947	di-2-butylamine;	(sec-C ₄ H ₉) ₂ NH;	[626-23-3]	0.04	0.0011
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