

THERMODYNAMICS

甲基-酰基-萘在庚烷、辛烷和十二烷中溶解度的测定与关联

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摘要 The solubilities of 2-methyl-6-acetylnaphthalene (2,6-MAN) and 2-methyl-7-acetylnaphthalene (2,7-MAN) in n-heptane, n-octane, and n-dodecane were measured, respectively, from 273.15 to 319.15K using an analytical method. On the basis of the thermodynamics theory of solid-liquid equilibrium, a model was derived to relate the solubilities with temperature. Using the least square method, the parameters of the model, the fusion enthalpies $\Delta_{fus}H$ and the Margules equation coefficients A12 and A21 of 2,6-MAN and 2,7-MAN in n-heptane, n-octane, and n-dodecane—were obtained by regressing the experimental data. The average deviation of the model was 1.70%.

关键词 [2-methyl-6-acetylnaphthalene](#), [2-methyl-7-acetylnaphthalene](#), [solubility](#), [analytical method](#)

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Experimental measurement and correlation of the solubility of methyl-acyl-naphthalene in n-heptane, n-octane, and n-dodecane

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Abstract The solubilities of 2-methyl-6-acetylnaphthalene (2,6-MAN) and 2-methyl-7-acetylnaphthalene (2,7-MAN) in n-heptane, n-octane, and n-dodecane were measured, respectively, from 273.15 to 319.15K using an analytical method. On the basis of the thermodynamics theory of solid-liquid equilibrium, a model was derived to relate the solubilities with temperature. Using the least square method, the parameters of the model, the fusion enthalpies $\Delta_{fus}H$ and the Margules equation coefficients A12 and A21 of 2,6-MAN and 2,7-MAN in n-heptane, n-octane, and n-dodecane—were obtained by regressing the experimental data. The average deviation of the model was 1.70%.

Key words [2-methyl-6-acetylnaphthalene](#); [2-methyl-7-acetylnaphthalene](#); [solubility](#); [analytical method](#)

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