

عنوان مقاله:

Effect of Temperature Functions $a(T)$ on Prediction of CH₄ and N₂ Gas Hydrate Formation Conditions

محل انتشار:

هفتمین کنگره ملی مهندسی شیمی (سال: 1390)

تعداد صفحات اصل مقاله: 7

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خلاصه مقاله:

In this research, by using Van der Waals–Plauteeuw model for solid hydrate phase and equations of state with mixing rules ($j-j$ model) for calculation of fugacity of components in gas and liquid phases, three phase equilibrium conditions (VLH)W have been predicted in two hydrate mixtures contain binary systems (CH₄-H₂O) and (N₂-H₂O). The pressure of hydrate formation has been calculated with four EOSs, VPT, PR, SRK and SW by using Danesh mixing rule. Also three temperature functions $a(T)$ in PR equation of state have been studied that contain Soave, Tribble–Bishnoi (TB) and Nasrifar-Moshfeghiyan (NM). Mixing rule interaction parameters in each mixture have been optimized by () W VL data and then optimized parameters have been used for three phase equilibrium calculations (VLH)W. Results show, for ($T > 302$ K and $P > 775$ bar) PR equation of state and for ($T < 290$ K and $P < 159$ bar) SRK equation of state have more accuracy in methane hydrate. In nitrogen hydrate mixture, SRK equation of state with 3% has minimum of error. Also TB temperature function with 7% error for (CH₄-H₂O) system and NM function with 8% error for (N₂-H₂O) system in PR equation of state have lower error than others functions.

کلمات کلیدی:

Methane hydrate, Nitrogen hydrate, Equation of state, Gas hydrate, Three Phase equilibrium

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