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Y Coulier, A Lowe, Jean Yves Coxam, Karine Ballerat-Busserolles. Liquid-liquid equilibriums in aqueous solutions of demixing amines loaded with gas for CO 2 capture processes. ECCE10+ECAB3+EPIC5, Sep 2015, NICE, France. hal-01226161

HAL Id: hal-01226161 https://hal.science/hal-01226161v1

Submitted on 8 Nov 2015

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Liquid-liquid equilibriums in aqueous solutions of demixing amines loaded with gas for CO₂ capture processes

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Introduction



Aqueous solutions of amines are selective carbon dioxide (CO_2) absorbents [1,2]. They are used in CO_2 capture processes working on absorption/desorption cycles.

Demixing amines may be valuable for new capture processes [3]. These amines demonstrate a liquid-liquid phase separation [4] which can be used to reduce energy costs of desorption (absorbent regeneration).

Aim of this paper : presentation of a method developed in the laboratory to study precisely the LLE in the binary solutions water + amines and the ternary mixtures water+amine+CO₂, as a function of the pressure and the CO₂ loading charge.

Amines : N-methylpiperidine





Institut de Chimie de

Clermont-Ferrand







Below temperature of demixion Above temperature of demixion

Sapphire cell **Hastelloy cell** 263 – 423 K Temperature T Room T – 428 K Control of T Thermostatic bath Heat tape 0.1 – 40 MPa 0.1 – 8 MPa Pressure *p* Control of p **Buffer volume** Buffer volume 15 to 20 mL Volume of the cell 5 mL adjustable Visualization of the Through the Full sample sapphire window sample

Below temperature of demixion Above temperature of demixion

Results and Conclusions

Constant loading charge

Temperatures of liquid-liquid phase separation for solutions of water + N-methylpiperidine (NMP) + CO_2 (figure 1) and for solutions of water + 2-methylpiperidine (2MP) + CO_2 (figure 2) versus composition of the binary mixture water + methylpiperidine at different loading charges α . For 2MP, data at α = 0 are from reference [5]





Constant amine-water composition

Temperatures of liquid-liquid phase separation for solutions of water + N-methylpiperidine (NMP) + CO_2 (figure 3) and for solutions of water + 2-methylpiperidine (2MP) + CO_2 (figure 4) versus loading charge α at different compositions of the binary mixture water + methylpiperidine. Open symbols : Thar Inst. Cell; full symbol : Sapphire cell



LC31 - 303 R

α = 0.33



- Good agreement between the two methods

2-methylpiperidine is a better candidate.

- LCST is strongly influenced by the presence of the dissolved gas
- N-methylpiperidine cannot be used in a process at high concentration (40%wt) as the CO₂ drastically lowered the LCST;

- Measurements were realized at 5 bar. Ternary mixtures are prepared below the temperature of liquid-liquid phase separation.

- <u>Uncertainties</u>

- composition : $\int_{\alpha}^{\alpha} = \pm 0.01$

	Water + amine	Water + amine + CO_2 (α = 0.2)	Water + amine + CO_2 (x = 0.05)
-methylpiperidine	LCST = 341 K x = 0.07	LCST = 343 K x = 0.04	
	1 CCT - 210 V		1 CCT - 202 V

- 2TO V

x = 0.07

 $\int x : \pm 0.001$

- Pressure : ± 0.1 bar , - Temperature : ± 0.3 K

ANR – NSERC joint program DACOOTA

N-methylpiperidine

The objective of the project, co-supported by ANR in France and NSERC in Canada, is to analyze the structure-properties relationships for different substituted piperidines.

Thermodynamic properties (enthalpies, heat capacities, volumes, phase equilibriums) are measured in Clermont-Ferrand (France). [6] Speciation in solution using Raman spectroscopy at high temperature and pressure are determined in Guelph (canada).

Modeling using activity coefficient models and molecular simulation will complete the program.

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