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SCALING DOWN OF FLASH CALCULATIONS OF NATURAL GAS
WITH THE AID OF PSEUDOCOMPONENTS

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A Thesis Submitted in Partial Fulfillment of the Requirements
for the Degree of Master of Engineering

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Thesis Title Scaling Down of Flash Calculations of Natural Gas with
the Aid of Pseudocomponents

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หัวขอวิทยานิพนธ์	การลดปริมาณการคำนวนแพลชของกําชธรรมชาติโดยอาศัยองค์ประกอบตัวแทน
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บทคัดย่อ

การคำนวณสมดุลย์ไอ-ของเหลวของกําชธรรมชาติจะเกี่ยวข้องกับการทำความสะอาดที่สมดุลย์ด้วยสมการสภาวะซึ่งในที่นี้เลือกใช้สมการสภาวะของ Soave-Redlich-Kwong และการทำความสะอาดประกอบในเฟสไอและของเหลวโดยวิธีแพลช การคำนวนมักจะเป็นการคำนวนแบบ iterative หรือ trial and error จึงต้องใช้คอมพิวเตอร์ช่วยในการคำนวน เมื่อจากองค์ประกอบในกําชธรรมชาติมีมากทำให้ลืมเปลี่ยนเวลาการคำนวน เพื่อเป็นการประหยัดเวลาการคำนวนจึงได้เสนอแบบจำลองที่อาศัยองค์ประกอบตัวแทนมาช่วย แบบจำลองนี้สามารถลดเวลาการคำนวนลงได้ 32% ของเวลาการคำนวนแบบธรรมดากลางๆ และมีเบอร์เชนต์ความผิดพลาดอยู่ในช่วง 0.6 ถึง 3.1 แบบจำลองนี้ใช้กับระบบความดันต่ำและปานกลาง สำหรับช่วงความดันและอุณหภูมิที่ทำการศึกษาคือ 150.0 psia ถึง 1115.0 psia และ -8.0°F ถึง 50.0°F

ศูนย์วิทยทรัพยากร
จุฬาลงกรณ์มหาวิทยาลัย

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ABSTRACT

The prediction of vapor-liquid equilibrium of natural gas deals with the calculation equilibrium K values based on the Soave-Redlich-Kwong equation of state and the phase compositions by flash calculations. Service of a computer is required due to the iterative nature of the calculation. Because of many equations to be solved simultaneously, much computing time is required. Therefore, in this work, the pseudocomponents model was introduced to scale down the computing time. With this model, the computing time was saved to 32% of the ordinary method. The percentage average deviation was the range of 0.6 to 3.1. This proposed model was tested at low and moderate pressure ranging from 150.0 psia to 1115.0 psia and a temperature range of -8.0°F to 50.0°F .



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TABLE OF CONTENTS

	Page
ABSTRACT IN THAI.....	iii
ABSTRACT IN ENGLISH.....	iv
ACKNOWLEDGEMENT.....	v
TABLE OF CONTENTS.....	vi
LIST OF TABLES.....	vii
LIST OF FIGURES.....	xvi
NOMENCLATURES.....	xviii
 CHAPTER	
1. INTRODUCTION.....	1
1.1 Introduction.....	1
1.2 Objective.....	2
1.3 Scope.....	2
2. VAPOR-LIQUID EQUILIBRIUM	3
2.1 Phase Behavior and Physical Properties of Natural Gas.....	3
2.2 Thermodynamics of Vapor-Liquid Equilibrium.....	9
2.3 Equations of State.....	10
3. VAPOR-LIQUID EQUILIBRIUM CALCULATIONS.....	18
3.1 Flash Calculation.....	18
3.2 Bubble and Dew Point Calculations.....	21
3.2.1 Bubble Point Calculation.....	22
3.2.2 Dew Point Calculation.....	22

	Page
3.2.2 Dew Point Calculation	22
3.2.3 Checking Phase Conditions for a Mixture	23
3.3 Method of Calculation	24
3.3.1 Newton-Raphson Method	25
3.3.2 Desirable Characteristics	27
3.3.3 Initial Estimate and Tolerance	28
4. PROPOSED GENERAL MODEL WITH PSEUDOCOMPONENTS	
4.1 Rationale	30
4.2 Basis	31
4.3 Algorithms	34
5. RESULTS AND DISCUSSION	
5.1 The Soave-Redlich-Kwong Equation of State	54
5.2 The Heavy-Pseudocomponent	54
5.3 Validity of Using Pseudocomponents	54
5.4 Study of Predicting the Compositions of Pseudocomponents	56
5.5 Comparison between the Ordinary Vapor-Liquid Equilibrium Calculation Model and the Proposed General Model with Pseudocomponents	58
6. CONCLUSION AND RECOMMENDATION	119
BIBLIOGRAPHIES	120
APPENDIX	125
BIOGRAPHY	147

LIST OF TABLES

TABLE	Page
5.1 COMPARISON OF EQUILIBRIUM K VALUES FROM THE ORDINARY METHOD AND THE PROPOSED GENERAL MODEL FOR HEAVY-PSEUDOCOMPONENT (M = 6) AT T = 41.0°F P = 300.0 PSIA.....	60
5.2 COMPARISON OF EQUILIBRIUM K VALUES FROM THE ORDINARY METHOD AND THE PROPOSED GENERAL MODEL FOR HEAVY-PSEUDOCOMPONENT (M = 6) AT T = 30.0°F P = 300.0 PSIA.....	61
5.3 COMPARISON OF EQUILIBRIUM K VALUES FROM THE ORDINARY METHOD AND THE PROPOSED GENERAL MODEL FOR HEAVY-PSEUDOCOMPONENT (M = 6) AT T = 30.0°F P = 300.0 PSIA.....	62
5.4 COMPARISON OF EQUILIBRIUM K VALUES FROM THE ORDINARY METHOD AND THE PROPOSED GENERAL MODEL FOR HEAVY-PSEUDOCOMPONENT (M = 6) AT T = -8.0°F P = 1115.0 PSIA.....	63
5.5 COMPARISON OF EQUILIBRIUM K VALUES FROM THE ORDINARY METHOD AND THE PROPOSED GENERAL MODEL FOR HEAVY-PSEUDOCOMPONENT (M = 6) AT T = -10.0°F P = 800.0 PSIA.....	64
5.6 COMPARISON OF EQUILIBRIUM K VALUES FROM THE ORDINARY METHOD AND THE PROPOSED GENERAL MODEL FOR HEAVY-PSEUDOCOMPONENT (M = 6) AT T = 15.0°F P = 850.0 PSIA.....	65

TABLE	Page
5.7 COMPARISON OF EQUILIBRIUM K VALUES FROM THE ORDINARY METHOD AND THE PROPOSED GENERAL MODEL FOR HEAVY-PSEUDOCOMPONENT (M = 8) AT T = 41.0°F P = 300.0 PSIA.....	66
5.8 COMPARISON OF EQUILIBRIUM K VALUES FROM THE ORDINARY METHOD AND THE PROPOSED GENERAL MODEL FOR HEAVY-PSEUDOCOMPONENT (M = 8) AT T = 30.0°F P = 300.0 PSIA.....	67
5.9 COMPARISON OF EQUILIBRIUM K VALUES FROM THE ORDINARY METHOD AND THE PROPOSED GENERAL MODEL FOR HEAVY-PSEUDOCOMPONENT (M = 8) AT T = 30.0°F P = 300.0 PSIA.....	68
5.10 COMPARISON OF EQUILIBRIUM K VALUES FROM THE ORDINARY METHOD AND THE PROPOSED GENERAL MODEL FOR HEAVY-PSEUDOCOMPONENT (M = 8) AT T = -10.0°F P = 800.0 PSIA.....	69
5.11 COMPARISON OF EQUILIBRIUM K VALUES FROM THE ORDINARY METHOD AND THE PROPOSED GENERAL MODEL FOR HEAVY- PSEUDOCOMPONENT(M = 8) AT T = 15.0°F P = 850.0 PSIA.....	70
5.12 COMPARISON OF EQUILIBRIUM K VALUES FROM THE ORDINARY METHOD AND THE PROPOSED GENERAL MODEL FOR HEAVY-PSEUDOCOMPONENT (M = 8) AT T = -8.0°F P = 1115.0 PSIA.....	71

TABLE	Page
5.13 COMPARISON OF EQUILIBRIUM K VALUES FROM THE ORDINARY METHOD AND THE PROPOSED GENERAL MODEL FOR LIGHT-PSEUDOCOMPONENT (C_{1+}) AT $T = 41.0^{\circ}\text{F}$ $P = 300.0 \text{ PSIA}.....$	72
5.14 COMPARISON OF EQUILIBRIUM K VALUES FROM THE ORDINARY METHOD AND THE PROPOSED GENERAL MODEL FOR LIGHT-PSEUDOCOMPONENT (C_{1+}) AT $T = 30.0^{\circ}\text{F}$ $P = 300.0 \text{ PSIA}.....$	73
5.15 COMPARISON OF EQUILIBRIUM K VALUES FROM THE ORDINARY METHOD AND THE PROPOSED GENERAL MODEL FOR LIGHT-PSEUDOCOMPONENT (C_{1+}) AT $T = 30.0^{\circ}\text{F}$ $P = 300.0 \text{ PSIA}.....$	74
5.16 COMPARISON OF EQUILIBRIUM K VALUES FROM THE ORDINARY METHOD AND THE PROPOSED GENERAL MODEL FOR LIGHT-PSEUDOCOMPONENT (C_{1+}) AT $T = -10.0^{\circ}\text{F}$ $P = 800.0 \text{ PSIA}.....$	75
5.17 COMPARISON OF EQUILIBRIUM K VALUES FROM THE ORDINARY METHOD AND THE PROPOSED GENERAL MODEL FOR LIGHT-PSEUDOCOMPONENT (C_{1+}) AT $T = 15.0^{\circ}\text{F}$ $P = 850.0 \text{ PSIA}.....$	76
5.18 COMPARISON OF EQUILIBRIUM K VALUES FROM THE ORDINARY METHOD AND THE PROPOSED GENERAL MODEL FOR LIGHT-PSEUDOCOMPONENT (C_{1+}) AT $T = -8.0^{\circ}\text{F}$ $P = 1115.0 \text{ PSIA}.....$	77
5.19 COMPARISON OF THE INITIAL PHASE VALUES OF THE EXPERIMENT AND THE ESTIMATION VALUES FOR HEAVY- PSEUDOCOMPONENT AT $T = -8.0^{\circ}\text{F}$ $P = 1115.0 \text{ PSIA}....$	78

TABLE

	Page
5.20 COMPARISON OF THE INITIAL PHASE VALUES OF THE EXPERIMENT AND THE ESTIMATION VALUES FOR HEAVY- PSEUDOCOMPONENT AT T = 15.0° F P = 700.0 PSIA.....	79
5.21 COMPARISON OF THE INITIAL PHASE VALUES OF THE EXPERIMENT AND THE ESTIMATION VALUES FOR HEAVY- PSEUDOCOMPONENT AT T = 15.0° F P = 200.0 PSIA.....	80
5.22 COMPARISON OF THE INITIAL PHASE VALUES OF THE EXPERIMENT AND THE ESTIMATION VALUES FOR LIGHT- PSEUDOCOMPONENT AT T = 15.0° F P = 200.0 PSIA.....	83
5.23 COMPARISON OF THE INITIAL PHASE VALUES OF THE EXPERIMENT AND THE ESTIMATION VALUES FOR LIGHT- PSEUDOCOMPONENT AT T = 15.0° F P = 700.0 PSIA.....	84
5.24 COMPARISON OF THE INITIAL PHASE VALUES OF THE EXPERIMENT AND THE ESTIMATION VALUES FOR LIGHT- PSEUDOCOMPONENT AT T = -8.0° F P = 1115.0 PSIA.....	85
5.25 COMPARISON OF THE INITIAL PHASE VALUES OF THE EXPERIMENT AND THE ESTIMATION VALUES FOR IDEAL CASE AT T = 15.0° F P = 200.0 PSIA.....	88
5.26 COMPARISON OF THE INITIAL PHASE VALUES OF THE EXPERIMENT AND THE ESTIMATION VALUES FOR IDEAL CASE AT T = 15.0° F P = 700.0 PSIA.....	89
5.27 COMPARISON OF THE INITIAL PHASE VALUES OF THE EXPERIMENT AND THE ESTIMATION VALUES FOR IDEAL CASE AT T = -8.0° F P = 1115.0 PSIA.....	90
5.28 COMPARISON OF THE LIQUID PHASE AND THE VAPOR PHASE COMPOSITIONS FROM THE EXPERIMENT DATA AND THE ORDINARY VLE CALCULATION MODEL AT T = -10.0° F P = 500.0 PSIA..	91

TABLE	Page
5.29 COMPARISON OF THE LIQUID AND THE VAPOR PHASE COMPOSITIONS FROM THE EXPERIMENTAL DATA AND THE ORDINARY VLE CALCULATION MODEL AT T = 0.0° F P = 500.0 PSIA.....	92
5.30 COMPARISON OF THE LIQUID AND THE VAPOR PHASE COMPOSITIONS FROM THE EXPERIMENTAL DATA AND THE ORDINARY VLE CALCULATION MODEL AT T = 50.0° F P = 500.0 PSIA....	93
5.31 COMPARISON OF THE LIQUID AND THE VAPOR PHASE COMPOSITIONS FROM THE EXPERIMENTAL DATA AND THE ORDINARY VLE CALCULATION MODEL AT T = 47.0° F P = 815.0 PSIA....	94
5.32 COMPARISON OF THE LIQUID AND THE VAPOR PHASE COMPOSITIONS FROM THE EXPERIMENTAL DATA AND THE ORDINARY VLE CALCULATION MODEL AT T = -8.0° F P = 250.0 PSIA....	95
5.33 COMPARISON OF THE LIQUID AND THE VAPOR PHASE COMPOSITION FROM THE EXPERIMENTAL DATA AND THE ORDINARY VLE CALCULATION MODEL AT T = -8.0° F P = 1115.0 PSIA... .	96
5.34 COMPARISON OF THE LIQUID AND THE VAPOR PHASE COMPOSITIONS FROM THE EXPERIMENTAL DATA AND THE ORDINARY VLE CALCULATION MODEL AT T = -10.0° F P = 800.0 PSIA... .	97
5.35 COMPARISON OF THE LIQUID AND THE VAPOR PHASE COMPOSITIONS FROM THE EXPERIMENTAL DATA AND THE ORDINARY VLE CALCULATION MODEL AT T = 8.0° F P = 300.0 PSIA.....	98
5.36 COMPARISON OF THE LIQUID AND THE VAPOR PHASE COMPOSITIONS FROM THE EXPERIMENTAL DATA AND THE ORDINARY VLE CALCULATION MODEL AT T = 8.0° F P = 700.0 PSIA.....	99
5.37 COMPARISON OF THE LIQUID AND THE VAPOR PHASE COMPOSITIONS FROM THE EXPERIMENTAL DATA AND THE ORDINARY VLE CALCULATION MODEL AT T = 0.0° F P = 300.0 PSIA.....	100

TABLE	Page
5.38 COMPARISON OF THE LIQUID AND THE VAPOR PHASE COMPOSITIONS FROM THE EXPERIMENTAL DATA AND THE PROPOSED GENERAL MODEL AT $T = -10.0^{\circ}\text{F}$ $P = 500.0 \text{ PSIA}$	101
5.39 COMPARISON OF THE LIQUID AND THE VAPOR PHASE COMPOSITIONS FROM THE EXPERIMENTAL DATA AND THE PROPOSED GENERAL MODEL AT $T = 0.0^{\circ}\text{F}$ $P = 500.0 \text{ PSIA}$	102
5.40 COMPARISON OF THE LIQUID AND THE VAPOR PHASE COMPOSITIONS FROM THE EXPERIMENTAL DATA AND THE PROPOSED GENERAL DATA AT $T = 50.0^{\circ}\text{F}$ $P = 500.0 \text{ PSIA}$	103
5.41 COMPARISON OF THE LIQUID AND THE VAPOR PHASE COMPOSITIONS FROM THE EXPERIMENTAL DATA AND THE PROPOSED GENERAL MODEL AT $T = 47.0^{\circ}\text{F}$ $P = 815.0 \text{ PSIA}$	104
5.42 COMPARISON OF THE LIQUID AND THE VAPOR PHASE COMPOSITIONS FROM THE EXPERIMENTAL DATA AND THE PROPOSED GENERAL MODEL AT $T = -8.0^{\circ}\text{F}$ $P = 250.0 \text{ PSIA}$	105
5.43 COMPARISON OF THE LIQUID AND THE VAPOR PHASE COMPOSITIONS FROM THE EXPERIMENTAL DATA AND THE PROPOSED GENERAL MODEL AT $T = -8.0^{\circ}\text{F}$ $P = 1115.0 \text{ PSIA}$	106
5.44 COMPARISON OF THE LIQUID AND THE VAPOR PHASE AND COMPOSITIONS FROM THE EXPERIMENTAL DATA AND THE PROPOSED GENERAL MODEL AT $T = 0.0^{\circ}\text{F}$ $P = 300.0 \text{ PSIA}$	107
5.45 COMPARISON OF THE LIQUID AND THE VAPOR PHASE COMPOSITIONS FROM THE EXPERIMENTAL DATA AND THE PROPOSED GENERAL MODEL AT $T = 8.0^{\circ}\text{F}$ $P = 300.0 \text{ PSIA}$	108
5.46 COMPARISON OF THE LIQUID AND THE VAPOR PHASE COMPOSITIONS FROM THE EXPERIMENTAL DATA AND THE PROPOSED GENERAL MODEL AT $T = 8.0^{\circ}\text{F}$ $P = 700.0 \text{ PSIA}$	109

TABLE	Page
5.47 COMPARISON OF THE LIQUID AND THE VAPOR PHASE COMPOSITIONS FROM THE ORDINARY VLE CALCULATION MODEL AND THE PROPOSED GENERAL MODEL AT T = -10.0 °F P = 500.0 PSIA.....	110
5.48 COMPARISON OF THE LIQUID AND THE VAPOR PHASE COMPOSITIONS FROM THE ORDINARY VLE CALCULATION MODEL AND THE PROPOSED GENERAL MODEL AT T = 0.0 °F P = 500.0 PSIA.....	111
5.49 COMPARISON OF THE LIQUID AND THE VAPOR PHASE COMPOSITIONS FROM THE ORDINARY VLE CALCULATION MODEL AND THE PROPOSED GENERAL MODEL AT T = 50.0 °F P = 500.0 PSIA.....	112
5.50 COMPARISON OF THE LIQUID AND THE VAPOR PHASE COMPOSITIONS FROM THE ORDINARY VLE CALCULATION MODEL AND THE PROPOSED GENERAL MODEL AT T = 47.0 °F P = 815.0 PSIA.....	113
5.51 COMPARISON OF THE LIQUID AND THE VAPOR PHASE COMPOSITIONS FROM THE ORDINARY VLE CALCULATION MODEL AND THE PROPOSED GENERAL MODEL AT T = -8.0 °F P = 250.0 PSIA.....	114
5.52 COMPARISON OF THE LIQUID AND THE VAPOR PHASE COMPOSITIONS FROM THE ORDINARY VLE CALCULATION MODEL AND THE PROPOSED GENERAL MODEL AT T = -8.0 °F P = 1115.0 PSIA.....	115
5.53 COMPARISON OF THE LIQUID THE VAPOR PHASE COMPOSITIONS FROM THE ORDINARY VLE CALCULATION MODEL	

LIST OF FIGURE

FIGURE	Page
2.1 PHASE ENVELOPE FOR NATURAL GAS MIXTURE.....	4
2.2 POSSIBLE CRITICAL POINT LOCATIONS ALONG THE P-T ENVELOPE.....	7
2.3 RETROGRADE BEHAVIOR.....	8
2.4 TYPICAL BEHAVIOR.....	15
3.1 CONTINUOUS EQUILIBRIUM FLASH CALCULATION.....	19
3.2 NEWTON-RAPHSON METHOD.....	26
3.2a A CASE OF NO CONVERGENCE.....	26
3.2b A SECOND CASE OF NO CONVERGENCE.....	26
4.1 EFFECT OF TEMPERATURE AND MOLECULAR TYPE ON VAPOR-LIQUID EQUILIBRIUM K VALUES AT P 500 PSIA.....	32
4.2 EFFECT OF PRESSURE ON VAPOR-LIQUID EQUILIBRIUM AT 100 F.....	33
4.3 ALGORITHM OF THE ORDINARY VAPOR-LIQUID EQUILIBRIUM CALCULATION MODEL.....	38
4.4 ALGORITHM OF THE PROPOSED GENERAL MODEL WITH PSEUDOCOMPONENTS.....	39
4.5 ALGORITHM FOR GROUPING LIGHT-PSEUDOCOMPONENT (SUBROUTINE LIGHT).....	43
4.6 ALGORITHM FOR GROUPING HEAVY-PSEUDOCOMPONENT (SUBROUTINE HEAVY).....	44
4.7 ALGORITHM FOR K VALUES DETERMINATION FOR THE SOAVE- REDLICH-KWONG EQUATION OF STATE(SUBROUTINE SRK).....	45

TABLE	Page
AND THE PROPOSED GENERAL MODEL AT T = 8.0°F	
·P = 300.0 PSIA.....	116
5.54 COMPARISON OF THE LIQUID AND THE VAPOR PHASE	
COMPOSITIONS FROM THE ORDINARY VLE CALCULATION MODEL	
AND THE PROPOSED GENERAL MODEL AT T = 8.0°F	
· P = 200.0 PSIA.....	117
5.55 COMPARISON OF THE LIQUID AND THE VAPOR PHASE	
COMPOSITIONS FROM THE ORDINARY VLE CALCULATION MODEL	
AND THE PROPOSED GENERAL MODEL AT T = 8.0°F	
P = 700.0 PSIA.....	118

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จุฬาลงกรณ์มหาวิทยาลัย

FIGURE	Page
4.8 ALGORITHM FOR THE SOLUTION OF THE CUBIC EQUATION OF (Z-COMPRESSIBILITY) (SUBROUTINE ROOTZ).....	47
4.9 ALGORITHM OF EQUILIBRIUM FLAS CALCULATION (SUBROUTINE FLASH).....	48
4.10 ALGORITHM FOR VAPOR PRESSURE CALCULATION.....	49
4.11 ALGORITHM OF BUBBLE POINT TEMPERATURE CALCULATION (SUBROUTINE DEWBUB).....	50
4.12 ALGORITHM OF DEW POINT TEMPERATURE CALCULATION (SUBROUTINE DEWBUB).....	51
4.13 ALGORITHM OF BUBBLE POINT PRESSURE CALCULATION (SUBROUTINE DEWBUB).....	52
4.14 ALGORITHM OF DEW POINT PRESSURE CALCULATION (SUBROUTINE DEWBUB).....	53
5.1 DEFINITION OF AREAS FOR ERROR ANALYSIS.....	56
5.2 CRITICAL LOCI OF CARBON DIOXIDE-PARAFFIN SYSTEMS.....	56
5.3 THE COMPARISON OF EXPERIMENTAL VALUES (ABCISSA) AND ESTIMATED VALUES (ORDINATE) IN THE LIQUID PHASE OF HEAVY-PSEUDOCOMPONENTS.....	81
5.4 THE COMPARISON OF EXPERIMENTAL VALUES (ABCISSA) AND ESTIMATED VALUES (ORDINATE) IN THE VAPOR PHASE OF HEAVY-PSEUDOCOMPONENTS.....	82
5.5 THE COMPARISON OF EXPERIMENTAL VALUES (ABCISSA) AND ESTIMATED VALUES (ORDINATE) IN THE LIQUID PHASE OF LIGHT-PSEUDOCOMPONENTS.....	86
5.6 THE COMPARISON OF EXPERIMENTAL VALUES (ABCISSA) AND ESTIMATED VALUES (ORDINATE) IN THE VAPOR PHASE OF LIGHT-PSEUDOCOMPONENTS.....	87

NOMENCLATURES



- A : Parameter in the SRK equation of state
- a_{c_i} : Parameter in the SRK equation of state
- ($a_{c\alpha}$) : Parameter in the SRK equation of state
- B : Parameter in the SRK equation of state
- b : Parameter in the SRK equation of state
- F : Degree of Freedom
- f_i : Fugacity of component i
- G : Gibb's Free Energy
- K_i : Equilibrium K values of component i
- k_{ij} : Binary interaction parameter
- L : Liquid flow rate, Liquid phase
- M : Subscript for the lightest component in the heavy pseudocomponent
- m_i : Parameter of the SRK equation of state
- N : The total number of component
- P : Total pressure
- P_i : Vapor pressure (saturation pressure of a pure species)
- P_{C_i} : Critical pressure of component i
- P_{R_i} : Reduced pressure = P/P_{C_i}
- Q : Parameter in the SRK equation of state
- R : Unibersal gasconstant
- T : Temperature
- T_{B_i} : The normal boiling point of component i
- T_{C_i} : Critical temperature of component i
- T_{R_i} : Reduced temperature = T/T_{C_i}

- x_1 : The liquid phase composition, the first approximation of root
- y_1 : The vapor phase composition
- z : Compressibility factor
- z_1 : Feed composition

GREEK LETTERS

- α : Parameter in The SRK equation of state
- β : Parameter in BWRS equation of state
- γ : Parameter in BWRS equation of state
- ρ : Density
- Δ : Difference operator
- ϵ : Convergerce tolerance
- μ : Chemical Potential
- Φ : Fugacity Coefficient of a component in a mixture
- ω : Acentric factor
- η : Parameter in the SRK equation of state

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