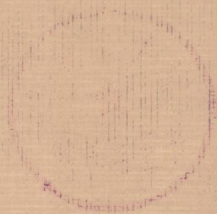


**PHYSICO-CHEMICAL PROPERTIES  
FOR CHEMICAL ENGINEERING**

**化学工学物性定数**

**Vol. 1**

**Edited by Physico-Chemical Properties Committee  
The Society of Chemical Engineers, Japan**



**MARUZEN CO., LTD., Tokyo**

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## PREFACE

The Japanese language edition of "Physical Properties of Matter" came to completion with the publishing of the tenth volume in 1970, however, due to popular request, it has been decided to publish additional works under the title of "Physico-Chemical Properties for Chemical Engineering." The main purpose of the new works is to introduce literatures on experimental data, experimental method of measurements, predictive method and theory of physico-chemical properties of matter related to equipment design in chemical engineering. Due to printing limitations, it is not possible to record the raw data. However, along with each of the chapter sections, an explanation of the means of measurement, margin of error, method of calculation, and remarks concerning the theory behind these will be included for whatever data can be measured. Great pains have been taken to make sure that the reader will be able to tell whether or not it is necessary to consult the original literature.

The chapter structure and the number of journals summarized are almost identical to those of "Physical Properties of Matter". But for brevity instructive reviews were not included in this publication.

Subsequent publication will be published every other year. One of the reasons for the use of English is that a friend in the U.S.A. suggested strongly that such valuable studies be published in a more widely used language.

At times it is intended to record material which is closely related to the subject under concern, but it may not be possible to go

## はしがき

日本語版「物性定数」は、一応1970年度の10巻をもって終刊したが、その後この種の書物を希望される多くの人の要望に答えるため、このたび「化学工学物性定数」の名のもとに本書を出版する運びとなった。本書の使命は、主として化学工業装置設計に関する物性定数のデータおよびその測定法、推算法およびその理論についての論文を広く紹介することである。紙面の都合から原データそのものの収録は不可能となったが、各章の諸項目それぞれについてデータの測定範囲、測定法、誤差、推算法とそれに関する理論の要点 (remarks) を記すことにより、読者が本書によって、原論文を取寄せる必要があるかないかの判定ができるように極力配慮したつもりである。

章の構成と摘録対象としてとりあげた雑誌数は、いままでの「物性定数」と比べてほとんど変わらない。書式を簡潔にするため、総説などは省略した。

本書は引続いて2年ごとに1集と同じ形式で発刊される。英語を用いた理由の一つは、米国の友人が、このような貴重な仕事はもう少し汎用性の多い用語を使用したらどうかと勧めてくれたことである。

時には関連の深い資料も収録する予定であるが、それでも本書に取上げている収録範囲は不十分であるかも知れない。既刊の数値集や便覧類とともに上手に活用して頂ければ、その便は非常に大きい

into sufficient depth. So, the use of other reference materials already on the market such as numerical tables and hand book is recommended for the sufficient use of this book.

For those who would like to refer to the original literature, or would like more detailed information than what is given in the bibliography, please refer to the notice for users (page 4) remarks.

Those who have started this work must be prepared to continue on in it as the field of study opens up along with the expansion of engineering. There's no doubt but that it is valuable work, but it takes a very generous person to take time off from his individual creative work, and devote it to such a non-profitable work as this is. The names of those who have so kindly given their cooperation are listed elsewhere. The fact that this publication has turned out so well is due greatly to the help of the first chairman Prof. Kazuo SATO, and the following chairman Prof. Mitsuho HIRATA, and also of the Editing Department of the Society of Chemical Engineers, Japan, and the Maruzen Co. Ltd. Publishing Department. And we owe typewriting service to Miss Etsuko NAKATA. We would like to express our deepest thanks to these people, and others who have been of help to us.

March 1977

Physico-Chemical Properties  
Committee Chairman  
Masahiro YORIZANE

ものと信ずる。

なお、原論文の参照を希望される方とか、「文献一覧表」よりさらに詳しい内容について知りたい方は凡例(4ページ)を参照されたい。

このような仕事は、いったんはじめた以上、学問の進展と工業の開発に伴って永久に続くものと覚悟しなければならぬであろう。きわめて大切な仕事であるには相違ないが、各自の創造的な仕事の他に利益を期待できないこのような仕事に時間をさくことは、きわめて奉仕的精神に富む人でなければできないことである。いままでこの仕事に協力して頂いた方々は別項に挙げたとおりである。

また、本書がこのように立派な形で刊行されるようになったのは、初代委員長の佐藤一雄教授およびその後を引き継がれた平田光穂教授の御尽力に負うところ最も大きく、そのほか化学工学協会編集課、および丸善株式会社出版部の諸氏の御努力のおかげである。また中田悦子女史には原稿のタイプをお願いした。これら多くの方々に対し、ここに深甚の謝意を表する次第である。

1977年3月

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頼 実 正 弘

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## 1. Structure of this book:

Each chapter comprises the following tables:

Table 1: English literature, data and equipment

Table 2: English literature, predictive method and theory

Table 3: Russian literature, data and equipment

Table 4: Russian literature, predictive method and theory

The parts (Part I, II, etc.) within each table are headed in such a way as to indicate the main points of the chapter.

## 2. Abbreviations and symbols:

Nomenclature will be explained before Table 1 and 3, if necessary.

Bibliographical notes will be placed by the following rule, and the bibliography itself will follow each individual chapter, without exception.

i) Chapters 1-8 will be indicated by symbols A-H.

ii) When the original literature is in Russian, an S will follow the chapter symbol.

iii) In each table, literature is shown by literature number with initial few letters of the name of author.

## 3. For those who want to refer to the original literature:

Send requests to...

Japan Information Center of Science and Technology

2-5, Nagata 2-chome, Chiyoda-ku Tokyo, 100, Japan

Make requests on...

JICST's stipulated request form, making sure to indicate the magazine number found on the right side of page five.

(This committee does not offer copying services.)

## 4. Concerning the services of this committee:

Those who want more detailed information than given in the bibliography should send a stamped self-addressed envelope to the address below. Replies will be made through the co-operation of the writers and reviewers mentioned elsewhere.

Physico-Chemical Properties Committee

The Society of Chemical Engineers, Japan

6-19, Kobinata 4-chome, Bunkyo-ku Tokyo, 112, Japan

## 1. 本文の構成

各章とも次の4つの表で構成されている。

表1: 原文が主として英文で、データ装置関係。

表2: 原文が主として英文で、推算法、理論に関するもの。

表3: 原文がロシア語で、内容は表1に同じ。

表4: 原文がロシア語で、内容は表2に同じ。

各表内の項 (Part I, II, etc.) は各章内の特長を生かした見出しがつけられている。

## 2. 略号と記号

Nomenclature は表1の前、または必要に応じて表3の前に記す。

文献記号は各章共通に下記のルールを用い、文献は各章末に添付してある。

i) 1~8章をA~Hの記号で代表させる。

ii) 原論文がロシア語のときは章記号の次にSをつける。

iii) 各表中で文献は、文献番号に著者の最初の数字を付して示されている。

## 3. 原論文の参照を希望される方へ

申込先...

日本科学技術情報センター (JICST)

〒100 東京都千代田区永田町2丁目5-2

申込方法...

JICST指定の申込用紙使用のこと。この場合雑誌記号には5ページ右端のJICST No. を明記のこと。

(当委員会では文献コピーサービスは致しません。)

## 4. 本委員会のサービスについて

文献一覧表以上のことをお知りになりたい場合は、原則として封書(返信用封書に宛名と切手同封のこと)で、下記へ連絡下さい。質問には、別項に挙げた執筆、摘録担当者が分担に応じて回答します。

宛先...

〒112 東京都文京区小日向4丁目6番19号

化学工学協会「物性定数委員会」

## INVESTICATED MAGAZINES AND THEIR JICST NUMBER

| No. | Full Name of Journal   | JICST No. |
|-----|--|-----------|
| 1   | Acta Chemica Scandinavica  | A 314A    |
| 2   | American Institute of Chemical Engineers Journal, New York                     | A 337A    |
| 3   | American Institute of Chemical Engineers Symposium Series                      | B 776A    |
| 4   | British Chemical Engineering <sup>see</sup> → Process Technology International | B 086A    |
| 5   | Bulletin of the Chemical Society of Japan                                      | G 450A    |
| 6   | Canadian Journal of Chemical Engineering                                       | B 227A    |
| 7   | Chemical Engineering   | B 251A    |
| 8   | Chemical Engineering Communications  | D 482A    |
| 9   | Chemical Engineering Journal   | D 723A    |
| 10  | Chemical Engineering Science   | B 254A    |
| 11  | Chemie-Ingenieur-Technik   | B 260A    |
| 12  | Chemical Reviews   | B 256A    |
| 13  | Collection of Czechoslovak Chemical Communications                             | B 288A    |
| 14  | Denki Kagaku oyobi Kogyo Butsuri Kagaku  | G 072A    |
| 15  | Doklady Akademij Nauk S.S.S.R.   | R 025A    |
| 16  | Tohoku Daigaku Hisuiyoeikagaku Kenkyusho Hokoku                                | F 342A    |
| 17  | Hydrocarbon Processing   | D 303A    |
| 18  | Industrial and Engineering Chemistry   | (A 531B)  |
| 19  | Izvestija Akademij Nauk S.S.S.R. (Seriya) Khimicheskaya                        | R 042A    |
| 20  | Journal of the American Chemical Society                                       | C 254A    |
| 21  | Journal of Applied Chemistry and Biotechnology                                 | C 264A    |
| 22  | Journal of Chemical and Engineering Data                                       | D 035B    |
| 23  | Journal of Chemical Engineering of Japan                                       | S 629A    |
| 24  | Journal of Chemical Physics  | C 275A    |
| 25  | Journal of the Chemical Society  | C 276A    |
| 26  | Journal of Chemical Thermodynamics   | D 634A    |
| 27  | Journal de Chimie Physique et de Physicochimie Biologique                      | C 277A    |
| 28  | Journal of the Indian Chemical Society   | C 296A    |
| 29  | Journal of Petroleum Technology  | C 330A    |
| 30  | Journal of Physical Chemistry  | C 334A    |
| 31  | Journal of Physical Chemistry of Liquid  | —         |
| 32  | Journal of Physical and Chemical Reference Data                                | —         |
| 33  | Journal of Research of the National Bureau of Standards. Sect A                | C 340B    |
| 34  | Kagaku Kōgaku  | F 099A    |
| 35  | Khimiya i tekhnologiya topliva i masel   | P 050A    |
| 36  | Khimicheskaja Promyshlennost   | R 049A    |
| 37  | Nippon Kagaku Kaishi   | F 226B    |
| 38  | Process Technology International   | B 086A    |
| 39  | Review of Physical Chemistry of Japan  | S 091A    |
| 40  | Transactions of the Faraday Society  | E 274A    |
| 41  | Transactions of the Institution of Chemical Engineers, London                  | E 282A    |
| 42  | Ukrainskii khimicheskii zhurnal  | R 203A    |
| 43  | Uspekhi Khimii   | P 117A    |
| 44  | Zeitschrift für Physikalische Chemie (Frankfurt)                               | E 443A    |
| 45  | Zeitschrift für Physikalische Chemie (Leipzig)                                 | E 444A    |
| 46  | Zhurnal Fizicheskoi Khimii   | R 136A    |
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CHAPTER 1 P-V-T RELATION AND LIQUID DENSITY

|                              |                               |
|------------------------------|-------------------------------|
| Nomenclature:                | (Greek letters)               |
| B; second virial coefficient | $\epsilon$ ; well depth       |
| C; third virial coefficient  | $\rho$ ; density              |
| c; molar concentration       | $\phi$ ; fugacity coefficient |
| d; molecular diameter        | $\omega$ ; acentric factor    |
| m; Stockmayer parameter      | (Subscripts)                  |
| n; refractive index          | c; critical                   |
| P; pressure                  | m; mixture                    |
| R; gas constant              | r; reduced                    |
| T; temperature               | s; saturated                  |
| V; molar volume              | (Superscripts)                |
| x; liquid mole fraction      | E; excess                     |
| Z; compressibility factor    | -; partial                    |

Table 1.1 Data

Part I PVT, Compressibility Factor, Expansion Factor, Virial Coefficient

| Substance         | Temperature<br>°C | Pressure<br>atm      | Remarks              | Lit.<br>(A) |
|-------------------|-------------------|----------------------|----------------------|-------------|
| AlCl <sub>3</sub> | 188~296           |                      | Saturated PVT        | 138         |
| CdCl <sub>2</sub> | 600               | 1 [bar]              | Z                    | 202         |
| CO <sub>2</sub>   | 0,16              | 30~100 [bar]         | Critical region      | 157         |
| Cs                | 704~1863          | 64~124               | PVT                  | 76E         |
| D <sub>2</sub> O  | 5~65              | 1~34 [bar]           | Z                    | 184         |
| H <sub>2</sub>    | 80.6 [°F]         | 2000~13000<br>[psia] | Z                    | 163<br>LOE  |
| He                | 70.0 [°F]         | 500~10000<br>[psia]  | Z                    | "           |
| Hg                | 30~150            | 8000 [bar]           | Thermal expansivity  | 97G         |
| H <sub>2</sub> O  | 25~150            | "                    | "                    | "           |
| H <sub>2</sub> O  | 5~35              | 100~300              | PVT                  | 250         |
| IF <sub>5</sub>   | 10~71             |                      | Z,B                  | 192         |
| KCl               | 800               | 1 [bar]              | Z                    | 202         |
| KNO <sub>3</sub>  | 400,500           | 1~5000 [bar]         | Z                    | "           |
| Kr                | -153~-53          | 5~3680               | Gas expansion method | 230         |
| Kr                | -128~-33          | 680                  | In dense fluid       | 231         |
| N <sub>2</sub>    | 80.6 [°F]         | 1500~13000<br>[psia] | Z                    | 163<br>LOE  |
| Ne                | -193~-143         | 2000                 | Z                    | 229         |

| Substance  | Temperature<br>°C      | Pressure<br>atm                 | Remarks                            | Lit.<br>(A) |
|--|------------------------|---------------------------------|------------------------------------|-------------|
| NH <sub>3</sub>  | -20~+40                | 1~1800                          | PVT                                | 149         |
| SF <sub>6</sub>  | -33~+46                | ~37.11                          | Critical region                    | 241         |
| Benzene  | 20~140                 |                                 | B                                  | 5AM         |
| 1-Bromopropane   | -75~+35                | 1~6000<br>[Kg/cm <sup>2</sup> ] | Z                                  | 127<br>JEN  |
| 1-Bromopentane   | -25~+125               | "                               | "                                  | "           |
| Butylaldehyde  | 30,40,50,60            | 5000~40000<br>[psig]            | Z                                  | 212<br>RUN  |
| CCl <sub>4</sub>   | 376.76~<br>1211.67[°R] | 0~600<br>[psia]                 | PVT                                | 178<br>MAR  |
| CCl <sub>4</sub>   | -19~+24                | 16[kbar]                        | PVT, Shock experiment              | 167         |
| Dichlorodifluoromethane  | 30~75                  |                                 | PVT, Z, B                          | 148         |
| Diethyl ether, Dibutyl-<br>ether, Diamyl ether,<br>Diethyl ether   | 20~35                  | ~1000<br>[psia]                 | Thermal pressure co-<br>efficients | 67<br>DRI   |
| Ethylene   | 50,100,150             |                                 | B                                  | 29B         |
| Ethylbenzene   | 10~90                  |                                 | Thermal expansivity                | 117         |
| Isopropanol  | 200~300                | 1000~8000<br>[psia]             | Z                                  | 245<br>TSE  |
| Methane  | 0~225                  | 20~600                          | PVT, Z                             | 193         |
| Methane  | -182~-28               | 350[bar]                        | PVT                                | 96G         |
| Methane  | 0~225                  | 1~350                           | Z                                  | 237         |
| Methyl ethyl ketone  | 20~60                  |                                 | Thermal expansivity                | 79F         |
| Octane   | -15~+85                | 1~1500                          | PVT                                | 20B         |
| Octane, Decane, Dodecane,<br>Tetradecane, Hexadecane   | -15~+80                | ~4000                           | PVT                                | 19<br>BEN   |
| Propane, Propene   | -20~+100<br>[°F]       | 21~223<br>[psia]                | Z                                  | 173<br>MAN  |
| n-Propanol   | 25~263.56              |                                 | Saturated PVT                      | 146         |
| Tetra-n-butylammonium<br>tetra-n-butylborate,<br>Tetraethylammonium tetra-<br>n-propylborate,<br>Tetra-n-propylammonium<br>tetraethylborate  | 90~160                 | ~5200                           | Z                                  | 98<br>GRI   |
| Ar+CO <sub>2</sub>   | 15                     | ~150                            | PVT, Z                             | 214         |
| Ar+KCl, Ar+KI, Ar+CaCl <sub>2</sub> , H <sub>2</sub> +<br>KOH, O <sub>2</sub> +KOH, Ethane+KCl,<br>Methane+KCl, Tetramethyl-<br>ammonium bromide+Ar,<br>Tetrabutylammonium<br>bromide+Ar, Tetramethyl- | 25                     |                                 | V̄                                 | 240<br>TIE  |

| Substance   | Temperature<br>°C     | Pressure<br>atm    | Remarks                         | Lit.<br>(A)       |
|---|-----------------------|--------------------|---------------------------------|-------------------|
| ammonium bromide+Methane,<br>Tetrabutylammonium<br>bromide+Methane  | 25                    |                    | $\bar{V}$                       | 240<br>TIE        |
| CO <sub>2</sub> +H <sub>2</sub> O, N <sub>2</sub> O+H <sub>2</sub> O,<br>Ethane+H <sub>2</sub> O, Ar+H <sub>2</sub> O   | 25, 50, 75,<br>100    | 22.4~50.8          | Cross B from<br>solubility data | 55<br>COA         |
| KNO <sub>3</sub> +Ca(NO <sub>3</sub> ) <sub>2</sub>   | 150~200               | 0~1500 [bar]       | Z                               | 202               |
| N <sub>2</sub> +H <sub>2</sub> , Methane+Ar+N <sub>2</sub> +H <sub>2</sub>  | -173~+17              | 150                | Joule-Thomson coeff.            | 93G               |
| N <sub>2</sub> , He, He+N <sub>2</sub>  | 38~177                | 100~5000<br>[psia] | Z                               | 179<br>MAT        |
| PbBr <sub>2</sub> +KBr, PbBr <sub>2</sub> +CsBr   | 1055~1164             | 22~43 [mmHg]       | Gas PVT                         | 243               |
| Tetrachloroformammonium-<br>bromide+H <sub>2</sub> O,<br>Tetramethylammonium-<br>bromide+H <sub>2</sub> O,<br>Tetraisopropylammonium-<br>bromide+H <sub>2</sub> O | 5.7, 24.2             |                    | Thermal expansivity             | 166<br>LUC        |
| Ar+CF <sub>4</sub>  | 100                   |                    | Cross virial coeff.             | 220               |
| Ar+Neopentane   | 50                    | 10~247             | Saturated PVT                   | 211               |
| CO <sub>2</sub> +N <sub>2</sub> , Methane+CO <sub>2</sub>   | -20, 0, 20            | 20~150             | PVT, Z                          | 9AR               |
| Difluorochloromethane+H <sub>2</sub> O  | 49.9~151.8            | 1~1800             | B                               | 235               |
| Ethane+H <sub>2</sub> S   | 50~125                | 5000 [psia]        | Z                               | 242               |
| Ethylene+Ar   | 50                    |                    | B                               | 29B               |
| Ethylene+NH <sub>3</sub>  | 30, 41                |                    | B                               | 208               |
| Methanol+N <sub>2</sub> , Methanol+Ar,<br>Methanol+Methane, Methanol<br>+Ethylene, Methanol<br>+Ethane, Methanol+CO <sub>2</sub> ,<br>Methanol+H <sub>2</sub> O   | 15, 25, 37,<br>50, 60 |                    | Cross B from<br>solubility data | 112<br>HEM        |
| Pyridine Hydrochloride<br>+ZnCl <sub>2</sub>  | 120~170               | 0~1500<br>[bar]    | Z                               | 202<br>POL        |
| SF <sub>6</sub> +CF <sub>4</sub>  | -1.5~+150             | 0~27               | Z, B, C                         | 221               |
| SF <sub>6</sub> +CF <sub>4</sub> , SF <sub>6</sub> +CF <sub>4</sub>   | "                     | "                  | "                               | "                 |
| Benzene+Cyclohexane   | 25                    |                    | Z                               | 77E               |
| Benzene+Isopropanol+H <sub>2</sub> O  | -10~+80               | 1~2800 [bar]       | PVT                             | 115               |
| 1-Chlorooctane, 1-Chloro-<br>decane, Capronitrile,<br>Caprylonitrile, Capri-<br>nitrile, Lauronitrile   | 16~153                |                    | PVT                             | 181<br>MEY        |
| Ethane+Methane  | -59~+35 [°F]          | 50~950<br>[psia]   | Joule-Thomson coeff.            | 4                 |
| Ethylene+Methanol   | 31, 60                |                    | B                               | ALK<br>208<br>RAT |

| Substance  | Temperature<br>°C | Pressure<br>atm | Remarks                     | Lit.<br>(A) |
|--|-------------------|-----------------|-----------------------------|-------------|
| Formic acid+H <sub>2</sub> O, Acetic acid+H <sub>2</sub> O, Propionic acid+H <sub>2</sub> O, Isobutyric acid+H <sub>2</sub> O                | 25~55             | 500~2500        | Z                           | 143<br>KOR  |
| Glycine+H <sub>2</sub> O, Glycolamide+H <sub>2</sub> O, Alanine+H <sub>2</sub> O, Lactamide+H <sub>2</sub> O, Glycylglycine+H <sub>2</sub> O | 25                | ~1000           | PVT                         | 254<br>YAY  |
| Hexane+Diethylether  | 39,60             |                 | B                           | 208         |
| Hexane+Methanol  | 50                |                 | Z, B                        | 62D         |
| Hexane+Propane   | 40~220            | 800 [psia]      | PVT                         | 132         |
| Methane+CF <sub>4</sub> , Butane+C <sub>4</sub> F <sub>10</sub>  | 25, 50, 100       |                 | Cross virial coeff.         | 220         |
| Ethane+CF <sub>4</sub> , Butane+CF <sub>4</sub> , Hexane+CF <sub>4</sub> , Methane+C <sub>2</sub> F <sub>6</sub>                             | 50                |                 | "                           | 220<br>SIE  |
| Ethane+C <sub>2</sub> F <sub>6</sub>   | 50, 100           |                 | "                           | "           |
| Butane+C <sub>2</sub> F <sub>6</sub> , Pentane+C <sub>2</sub> F <sub>6</sub>   | 50                |                 | "                           | "           |
| Propane+C <sub>3</sub> F <sub>8</sub>  | 50, 100           |                 | "                           | "           |
| Methane+C <sub>4</sub> F <sub>10</sub> , Ethane+C <sub>4</sub> F <sub>10</sub> , Hexane+C <sub>4</sub> F <sub>10</sub>                       | 50                |                 | "                           | "           |
| Methane+C <sub>5</sub> F <sub>12</sub>   | 100               |                 | "                           | "           |
| Pentane+C <sub>5</sub> F <sub>12</sub>   | 50, 100           |                 | "                           | "           |
| Methane+C <sub>6</sub> F <sub>14</sub>   | 100               |                 | "                           | "           |
| Hexane+C <sub>6</sub> F <sub>14</sub>  | 50, 100           |                 | "                           | "           |
| Neopentane+CF <sub>4</sub> , CF <sub>4</sub> +SF <sub>6</sub>  | 50                |                 | "                           | "           |
| Methane+Neopentane   | 25                | 11~149          | Saturated PVT               | 211         |
| Neopentane, bicyclo[2,2,2]-Octane, Adamantane  | 25                |                 | Z                           | 33<br>BRE   |
| Pentane+Methanol   | 31, 40            |                 | B                           | 208         |
| p-Xylene+Benzene, p-Dioxane+Benzene  | 25, 40            |                 | Z <sup>E</sup>              | 135<br>KHA  |
| Binary gas-gas mixtures  |                   |                 | Review, V <sup>E</sup> data | 17<br>BAT   |

## Part II Liquid Density

|                                       |             |              |                  |     |
|---------------------------------------|-------------|--------------|------------------|-----|
| AgNO <sub>3</sub> , TlNO <sub>3</sub> | 433~583 [K] |              | Density equation | 125 |
| AlCl <sub>3</sub>                     | 188~296     |              | Saturated values | 138 |
| AlCl <sub>3</sub>                     | 180~360     |              | Density equation | 219 |
| N <sub>2</sub>                        | 182~-173    |              |                  | 162 |
| ND <sub>3</sub>                       | -25~+30     |              |                  | 3   |
| Fluorobenzene                         | 30, 60, 100 | 1~2000 [bar] |                  | ALE |
| NiCl <sub>2</sub>                     | 1015~1058   |              |                  | 84G |
| Pb                                    | 344~652     |              | Density equation | 217 |
| XeF <sub>2</sub>                      | 90~115      |              | Density equation | 191 |
| XeF <sub>4</sub>                      | 50~110      |              | "                | "   |

| Substance   | Temperature<br>°C | Pressure<br>atm | Remarks          | Lit.<br>(A) |
|---|-------------------|-----------------|------------------|-------------|
| Acetic acid   | 20, 30, 40        |                 |                  | 16B         |
| Benzene   | 20~289            | 0.1~48.0        |                  | 5AM         |
| Benzene, Toluene, o-Xylene,<br>m-Xylene, p-Xylene, Ethyl<br>benzene, Isopropyl benzene,<br>Naphthalene, 2-Methyl<br>naphthalene   | 20~217            |                 | Density equation | 104<br>HAL  |
| CF <sub>3</sub> PCl <sub>2</sub> , CF <sub>3</sub> PBr <sub>2</sub>   | 20~69             |                 |                  | 131         |
| 1,2-Dichloroethane  | 35~65             |                 |                  | 47C         |
| 1,4-Dichlorobutane  | "                 |                 |                  | "           |
| 1,6-Dichlorohexane  | "                 |                 |                  | "           |
| 1,8-Dichlorooctane  | "                 |                 |                  | "           |
| 1,10-Dichlorodecane   | "                 |                 |                  | "           |
| 2,3-Dimethyl-2-butene,<br>3,3-Dimethyl-1-butene   | 20~50             |                 | Density equation | 15<br>BAG   |
| Dimethyl-3,5-pyridine   | 20, 25, 35, 45    |                 |                  | 182         |
| Dimethyl-2,6-pyridine   | "                 |                 |                  | "           |
| Ethylacetate,<br>Isoamyl alcohol,<br>Benzyl alcohol,<br>1,4-Dioxane   | 0~70              |                 | Density equation | 1<br>ABR    |
| Ethane, Ar+Kr, Ar+Xe  | 110.77 [K]        |                 |                  | 54C         |
| Ethane, Kr+Xe   | 161.36 [K]        |                 |                  | "           |
| Hexadecane, 1-Chlorobutane,<br>1-Chloropentane, 1-Chloro-<br>octane, 1-Chlorodecane,<br>1-Chlorohexane, 1-Chloro-<br>dodecane, 1-Chlorotetra-<br>decane, 1-Chlorohexadecane<br>, 1-Chlorooctadecane | 25                |                 |                  | 113<br>HER  |
| Hexafluoroacetone   |                   |                 | Tr=0.58~0.82     | 186         |
| Hexafluorobenzene   | 20, 25, 35, 45    |                 |                  | 182         |
| Methane   | -182~-153         |                 |                  | 162         |
| 2-Methyl-2,4-pentanediol,<br>Dipropylene glycol,<br>2-Ethyl-1,3-hexanediol,<br>Thiodiglycol   | 20~50             |                 |                  | 124<br>IKA  |
| Methyl-2-pyridine   | 20, 25, 35, 45    |                 |                  | 182         |
| Methyl-4-pyridine   | 20, 25, 35, 45    |                 |                  | "           |
| Octane  | -27~+122          | 1               |                  | 50C         |



| Substance  | Temperature<br>°C   | Pressure<br>atm | Remarks              | Lit.<br>(A) |
|--|---------------------|-----------------|----------------------|-------------|
| Octamethylcyclotetra-<br>siloxane  | 503.1~579.8<br>[K]  |                 |                      | 256<br>YOU  |
| cis-Pentene-2  | -129.72~+25         |                 |                      | 58C         |
| Pentane  | -151.39~+25         |                 |                      | "           |
| Polystyrene  | 20~110              |                 | Density equation     | 116         |
| Propionic acid, Butyric-<br>acid, Isobutyric acid,<br>Valeric acid   | 0, 10, 20, 30<br>40 |                 |                      | 16<br>BAR   |
| n-Propanol   | 25~263.56           |                 | Saturated properties | 146         |
| Pyridine   | 20, 25, 35, 45      |                 |                      | 182         |
| Tetrahydrofuran+H <sub>2</sub> O,<br>Tetrahydropyran+H <sub>2</sub> O,<br>2-Methyltetrahydrofuran+<br>H <sub>2</sub> O, 2,5-Dimethyltetra-<br>hydrofuran+H <sub>2</sub> O,<br>1,3-Dioxolane+H <sub>2</sub> O,<br>1,4-Dioxane+H <sub>2</sub> O,<br>Aziridine+H <sub>2</sub> O,<br>Azetidine+H <sub>2</sub> O,<br>Pyrrolidine+H <sub>2</sub> O,<br>Piperidine+H <sub>2</sub> O,<br>Hexamethylenimine+H <sub>2</sub> O,<br>Heptamethylenimine+H <sub>2</sub> O,<br>1-Methylpyrrolidine+H <sub>2</sub> O,<br>1-Methylpiperidine+H <sub>2</sub> O | 25                  |                 |                      | 37C         |
| Trifluoroacetonitrile  |                     |                 | Tr=0.61~0.92         | 186         |
| Trifluoroacetyl fluoride   | -93~-62             |                 | Density equation     | 194         |
| Trimethyl-2,4,6-pyridine   | 20, 25, 35, 45      |                 |                      | 182         |
| Ba+BaBr <sub>2</sub>   | 882~999             |                 |                      | 100         |
| Ba+BaCl <sub>2</sub>   | 986~1026            |                 |                      | "           |
| Bi+BiBr <sub>2</sub>   | 589 [K]             |                 | v <sup>E</sup>       | 234         |
| Bi+BiCl <sub>3</sub>   | 604 [K]             |                 | "                    | "           |
| Bi+BiI <sub>3</sub>  | 815 [K]             |                 | "                    | "           |
| CaCl <sub>2</sub> +MgCl <sub>2</sub> , SrCl <sub>2</sub> +MgCl <sub>2</sub> ,<br>BaCl <sub>2</sub> +MgCl <sub>2</sub>  | 900                 |                 |                      | 21<br>BER   |
| Catechol+Resorcinol  | 110, 115, 120       |                 |                      | 189         |
| Cd+CdCl <sub>2</sub> , Cd+CdBr <sub>2</sub>  | 1009 [K]            |                 | v <sup>E</sup>       | 234         |
| Cd+CdI <sub>2</sub>  | 793 [K]             |                 | "                    | "           |
| CsCl+BaCl <sub>2</sub>   | 965~1018            |                 |                      | 100         |
| CuCl <sub>2</sub> +HCl+H <sub>2</sub> O  | 30, 50, 70, 90      |                 |                      | 99G         |
| KCl+LaCl <sub>3</sub>  | 765~996             |                 | Density equation     | 53C         |