

# Formamide, N,N-dimethyl-

|                      |                                      |
|----------------------|--------------------------------------|
| Other names:         | DMF                                  |
|                      | DMF (amide)                          |
|                      | DMF (dimethylformamide)              |
|                      | DMFA                                 |
|                      | Dimethylamid kyseliny mravenci       |
|                      | Dimethylforamide                     |
|                      | Dimethylformamid                     |
|                      | Dimethylformamide                    |
|                      | Dimetilformamide                     |
|                      | Dimetylformamidu                     |
|                      | Dwumetyloformamid                    |
|                      | Formyldimethylamine                  |
|                      | HCON(CH3)2                           |
|                      | N,N-Dimethylformaldehyde             |
|                      | N,N-Dimethylformamide                |
|                      | N,N-Dimethylmethanamide              |
|                      | N-FORMYLDIMETHYLAMINE                |
|                      | NCI-C60913                           |
|                      | NSC-5356                             |
|                      | U-4224                               |
|                      | UN 2265                              |
|                      | dimethylformamide (DMF)              |
| Inchi:               | InChI=1S/C3H7NO/c1-4(2)3-5/h3H,1-2H3 |
| InchiKey:            | ZMXDDKWLCZADIW-UHFFFAOYSA-N          |
| Formula:             | C3H7NO                               |
| SMILES:              | CN(C)C=O                             |
| Mol. weight [g/mol]: | 73.09                                |
| CAS:                 | 68-12-2                              |

## Physical Properties

| Property code | Value           | Unit   | Source       |
|---------------|-----------------|--------|--------------|
| affp          | 887.50          | kJ/mol | NIST Webbook |
| basg          | 856.60          | kJ/mol | NIST Webbook |
| chl           | -1941.90        | kJ/mol | NIST Webbook |
| chl           | -1941.60 ± 1.20 | kJ/mol | NIST Webbook |

|         |                |        |  |
|---------|----------------|--------|--|
| dvisc   | 0.0008050      | Paxs   | Thermodynamic Properties of Salophen Schiff Base + Ionic Liquid ([Cnmlm][Br]) + Dimethylformamide Ternary Mixtures at 298.15 K   |
| dvisc   | 0.0008610      | Paxs   | Thermodynamic properties of ionic liquid, 1-hexyl-3-methylimidazolium bromide, + N-N'bis(2-pyridylmethylidene)-1,2-diiminoethane Schiff base + N,N-dimethylformamide solutions at T = (298.15 to 313.15) K |
| ea      | 0.01           | eV     | NIST Webbook   |
| gf      | -14.36         | kJ/mol | Joback Method  |
| hf      | -123.30        | kJ/mol | Joback Method  |
| hfl     | -239.40 ± 1.20 | kJ/mol | NIST Webbook   |
| hfl     | -239.00        | kJ/mol | NIST Webbook   |
| hfus    | 46.65          | kJ/mol | Enthalpies of vaporization of N,N-dialkyl monamides at 298.15K   |
| hfus    | 8.95           | kJ/mol | Thermodynamic properties of N,N-dimethylformamide and N,N-dimethylacetamide  |
| hvap    | 47.57          | kJ/mol | NIST Webbook   |
| hvap    | 46.90          | kJ/mol | NIST Webbook   |
| hvap    | 46.70 ± 0.50   | kJ/mol | NIST Webbook   |
| ie      | 9.12 ± 0.02    | eV     | NIST Webbook   |
| ie      | 9.25           | eV     | NIST Webbook   |
| ie      | 9.14           | eV     | NIST Webbook   |
| ie      | 9.45 ± 0.05    | eV     | NIST Webbook   |
| ie      | 9.13           | eV     | NIST Webbook   |
| ie      | 9.14           | eV     | NIST Webbook   |
| log10ws | 0.58           |        | Crippen Method   |
| logp    | -0.296         |        | Crippen Method   |
| mccvol  | 64.680         | ml/mol | McGowan Method   |
| nfpaf   | %!d(float64=2) |        | KDB  |
| nfpah   | %!d(float64=1) |        | KDB  |
| pc      | 4862.97        | kPa    | Joback Method  |
| rhoc    | 279.22 ± 5.85  | kg/m3  | NIST Webbook   |
| rinpol  | 752.00         |        | NIST Webbook   |
| rinpol  | 747.00         |        | NIST Webbook   |
| rinpol  | 750.00         |        | NIST Webbook   |
| rinpol  | 756.30         |        | NIST Webbook   |
| rinpol  | 745.00         |        | NIST Webbook   |
| rinpol  | 742.00         |        | NIST Webbook   |
| rinpol  | 790.00         |        | NIST Webbook   |

|        |               |   |  |
|--------|---------------|---|--|
| rinpol | 772.00        |   | NIST Webbook   |
| rinpol | 783.00        |   | NIST Webbook   |
| rinpol | 746.00        |   | NIST Webbook   |
| rinpol | 751.00        |   | NIST Webbook   |
| rinpol | 746.00        |   | NIST Webbook   |
| rinpol | 735.00        |   | NIST Webbook   |
| rinpol | 738.00        |   | NIST Webbook   |
| rinpol | 752.00        |   | NIST Webbook   |
| rinpol | 746.00        |   | NIST Webbook   |
| rinpol | 782.00        |   | NIST Webbook   |
| rinpol | 772.00        |   | NIST Webbook   |
| rinpol | 782.00        |   | NIST Webbook   |
| rinpol | 753.00        |   | NIST Webbook   |
| ripol  | 1276.00       |   | NIST Webbook   |
| ripol  | 1326.00       |   | NIST Webbook   |
| ripol  | 1361.00       |   | NIST Webbook   |
| ripol  | 1282.00       |   | NIST Webbook   |
| ripol  | 1312.00       |   | NIST Webbook   |
| ripol  | 1327.00       |   | NIST Webbook   |
| ripol  | 1290.00       |   | NIST Webbook   |
| ripol  | 1295.00       |   | NIST Webbook   |
| ripol  | 1276.00       |   | NIST Webbook   |
| ripol  | 1344.00       |   | NIST Webbook   |
| ripol  | 1361.00       |   | NIST Webbook   |
| ripol  | 1319.00       |   | NIST Webbook   |
| ripol  | 1328.00       |   | NIST Webbook   |
| ripol  | 1344.00       |   | NIST Webbook   |
| ripol  | 1333.00       |   | NIST Webbook   |
| ripol  | 1304.00       |   | NIST Webbook   |
| ripol  | 1325.00       |   | NIST Webbook   |
| ripol  | 1326.00       |   | NIST Webbook   |
| tb     | 426.00 ± 1.00 | K | NIST Webbook   |
| tb     | 426.00        | K | KDB  |
| tb     | 425.70        | K | Vapor liquid equilibria for water + acetic acid + (N,N-dimethylformamide or dimethyl sulfoxide) at 13.33 kPa   |
| tb     | 426.05        | K | Separation of azeotropic mixture (2, 2, 3, 3-Tetrafluoro-1-propanol + water) by extractive distillation: Entrainers selection and vapour-liquid equilibrium measurements |

|    |               |                      |   |
|----|---------------|----------------------|---|
| tb | 425.15        | K                    | Vapor Liquid Equilibrium Data for Methanol + tert-Butylamine + N,N-Dimethylformamide and Constituent Binary Systems at Atmospheric Pressure                   |
| tb | 428.15 ± 2.00 | K                    | NIST Webbook  |
| tb | 426.00 ± 3.00 | K                    | NIST Webbook  |
| tb | 425.85 ± 0.25 | K                    | NIST Webbook  |
| tb | 425.95 ± 0.50 | K                    | NIST Webbook  |
| tb | 426.05 ± 0.30 | K                    | NIST Webbook  |
| tb | 425.95 ± 0.70 | K                    | NIST Webbook  |
| tb | 426.20        | K                    | NIST Webbook  |
| tb | 426.21        | K                    | Isobaric Vapor Liquid Equilibria for the Binary and Ternary Systems of 2-Methyl-1-butanol, 2-Methyl-butanol Acetate, and Dimethylformamide (DMF) at 101.3 kPa |
| tc | 649.60        | K                    | KDB   |
| tc | 649.60 ± 0.60 | K                    | NIST Webbook  |
| tf | 212.72        | K                    | KDB   |
| tf | 212.70 ± 0.02 | K                    | NIST Webbook  |
| vc | 0.262         | m <sup>3</sup> /kmol | KDB   |

## Temperature Dependent Properties

| Property code | Value  | Unit    | Temperature [K] | Source        |
|---------------|--------|---------|-----------------|---------------|
| cpg           | 117.36 | J/mol×K | 386.37          | Joback Method |
| cpg           | 103.92 | J/mol×K | 329.14          | Joback Method |
| cpg           | 110.79 | J/mol×K | 357.75          | Joback Method |
| cpg           | 123.65 | J/mol×K | 414.98          | Joback Method |
| cpg           | 129.66 | J/mol×K | 443.59          | Joback Method |
| cpg           | 135.39 | J/mol×K | 472.20          | Joback Method |
| cpg           | 140.87 | J/mol×K | 500.82          | Joback Method |
| cpl           | 156.69 | J/mol×K | 298.00          | NIST Webbook  |
| cpl           | 146.05 | J/mol×K | 298.15          | NIST Webbook  |
| cpl           | 148.16 | J/mol×K | 298.15          | NIST Webbook  |
| cpl           | 149.28 | J/mol×K | 308.00          | NIST Webbook  |
| cpl           | 150.00 | J/mol×K | 298.15          | NIST Webbook  |
| cpl           | 148.36 | J/mol×K | 298.15          | NIST Webbook  |
| cpl           | 150.50 | J/mol×K | 298.15          | NIST Webbook  |
| cpl           | 150.80 | J/mol×K | 298.15          | NIST Webbook  |

|       |           |         |        |  |
|-------|-----------|---------|--------|--|
| cpl   | 120.50    | J/molxK | 298.00 | NIST Webbook   |
| cpl   | 148.00    | J/molxK | 298.15 | NIST Webbook   |
| cpl   | 148.00    | J/molxK | 298.00 | NIST Webbook   |
| cpl   | 150.80    | J/molxK | 298.15 | NIST Webbook   |
| cpl   | 146.00    | J/molxK | 298.15 | NIST Webbook   |
| cpl   | 152.00    | J/molxK | 298.15 | NIST Webbook   |
| dvisc | 0.0008045 | Paxs    | 298.15 | Viscosity of Urea<br>in the Mixture of<br>N,N-Dimethylformamide<br>and Water   |
| dvisc | 0.0008605 | Paxs    | 298.15 | Volumetric<br>properties of<br>ionic liquid<br>1,3-dimethylimidazolium<br>methyl sulfate +<br>molecular<br>solvents at T =<br>(298.15 - 328.15)<br>K |
| dvisc | 0.0007712 | Paxs    | 308.15 | Volumetric<br>properties of<br>ionic liquid<br>1,3-dimethylimidazolium<br>methyl sulfate +<br>molecular<br>solvents at T =<br>(298.15 - 328.15)<br>K |
| dvisc | 0.0006975 | Paxs    | 318.15 | Volumetric<br>properties of<br>ionic liquid<br>1,3-dimethylimidazolium<br>methyl sulfate +<br>molecular<br>solvents at T =<br>(298.15 - 328.15)<br>K |
| dvisc | 0.0006374 | Paxs    | 328.15 | Volumetric<br>properties of<br>ionic liquid<br>1,3-dimethylimidazolium<br>methyl sulfate +<br>molecular<br>solvents at T =<br>(298.15 - 328.15)<br>K |
| dvisc | 0.0007830 | Paxs    | 298.15 | Densities,<br>Viscosities, and<br>Conductivities of<br>Phosphonic Acid<br>Solutions in<br>N,N-Dimethylformamide<br>and Water                         |
| dvisc | 0.0006990 | Paxs    | 308.15 | Densities,<br>Viscosities, and<br>Conductivities of<br>Phosphonic Acid<br>Solutions in<br>N,N-Dimethylformamide<br>and Water                         |

|       |           |      |        |  |
|-------|-----------|------|--------|--|
| dvisc | 0.0006230 | Paxs | 318.15 | Densities, Viscosities, and Conductivities of Phosphonic Acid Solutions in N,N-Dimethylformamide and Water   |
| dvisc | 0.0005550 | Paxs | 328.15 | Densities, Viscosities, and Conductivities of Phosphonic Acid Solutions in N,N-Dimethylformamide and Water   |
| dvisc | 0.0008030 | Paxs | 298.15 | Excess Molar Volumes and Deviations in Viscosity of Binary Mixtures of N,N-Dimethylformamide with Aniline and Benzonitrile at (298.15, 303.15, 308.15, and 313.15) K |
| dvisc | 0.0007560 | Paxs | 303.15 | Excess Molar Volumes and Deviations in Viscosity of Binary Mixtures of N,N-Dimethylformamide with Aniline and Benzonitrile at (298.15, 303.15, 308.15, and 313.15) K |
| dvisc | 0.0007100 | Paxs | 308.15 | Excess Molar Volumes and Deviations in Viscosity of Binary Mixtures of N,N-Dimethylformamide with Aniline and Benzonitrile at (298.15, 303.15, 308.15, and 313.15) K |
| dvisc | 0.0006730 | Paxs | 313.15 | Excess Molar Volumes and Deviations in Viscosity of Binary Mixtures of N,N-Dimethylformamide with Aniline and Benzonitrile at (298.15, 303.15, 308.15, and 313.15) K |

|       |           |      |        |  |
|-------|-----------|------|--------|--|
| dvisc | 0.0008030 | Paxs | 298.15 | Densities, Viscosities, and Thermodynamic Properties of (N,N-Dimethylformamide + Benzene + Chlorobenzene) Ternary Mixtures at (298.15, 303.15, 308.15, and 313.15) K |
| dvisc | 0.0007560 | Paxs | 303.15 | Densities, Viscosities, and Thermodynamic Properties of (N,N-Dimethylformamide + Benzene + Chlorobenzene) Ternary Mixtures at (298.15, 303.15, 308.15, and 313.15) K |
| dvisc | 0.0007100 | Paxs | 308.15 | Densities, Viscosities, and Thermodynamic Properties of (N,N-Dimethylformamide + Benzene + Chlorobenzene) Ternary Mixtures at (298.15, 303.15, 308.15, and 313.15) K |
| dvisc | 0.0006730 | Paxs | 313.15 | Densities, Viscosities, and Thermodynamic Properties of (N,N-Dimethylformamide + Benzene + Chlorobenzene) Ternary Mixtures at (298.15, 303.15, 308.15, and 313.15) K |
| dvisc | 0.0008030 | Paxs | 298.15 | Densities and Viscosities of Binary Mixtures of N,N-Dimethylformamide with Benzyl Alcohol and Acetophenone at (298.15, 303.15, 308.15, and 313.15) K                 |

|       |           |      |        |  |
|-------|-----------|------|--------|--|
| dvisc | 0.0007560 | Paxs | 303.15 | Densities and Viscosities of Binary Mixtures of N,N-Dimethylformamide with Benzyl Alcohol and Acetophenone at (298.15, 303.15, 308.15, and 313.15) K |
| dvisc | 0.0007100 | Paxs | 308.15 | Densities and Viscosities of Binary Mixtures of N,N-Dimethylformamide with Benzyl Alcohol and Acetophenone at (298.15, 303.15, 308.15, and 313.15) K |
| dvisc | 0.0006730 | Paxs | 313.15 | Densities and Viscosities of Binary Mixtures of N,N-Dimethylformamide with Benzyl Alcohol and Acetophenone at (298.15, 303.15, 308.15, and 313.15) K |
| dvisc | 0.0008030 | Paxs | 298.15 | Density and Viscosity Studies of Binary Mixtures of N,N-Dimethylformamide with Toluene and Methyl Benzoate at (298.15, 303.15, 308.15, and 313.15) K |
| dvisc | 0.0007560 | Paxs | 303.15 | Density and Viscosity Studies of Binary Mixtures of N,N-Dimethylformamide with Toluene and Methyl Benzoate at (298.15, 303.15, 308.15, and 313.15) K |
| dvisc | 0.0007100 | Paxs | 308.15 | Density and Viscosity Studies of Binary Mixtures of N,N-Dimethylformamide with Toluene and Methyl Benzoate at (298.15, 303.15, 308.15, and 313.15) K |



|       |           |      |        |   |
|-------|-----------|------|--------|---|
| dvisc | 0.0006730 | Paxs | 313.15 | Density and Viscosity Studies of Binary Mixtures of N,N-Dimethylformamide with Toluene and Methyl Benzoate at (298.15, 303.15, 308.15, and 313.15) K  |
| dvisc | 0.0008586 | Paxs | 298.15 | Densities, Viscosities, and Sound Speeds of Some Acetate Salts in Binary Mixtures of Tetrahydrofuran and Methanol at (303.15, 313.15, and 323.15) K   |
| dvisc | 0.0007763 | Paxs | 308.15 | Densities, Viscosities, and Sound Speeds of Some Acetate Salts in Binary Mixtures of Tetrahydrofuran and Methanol at (303.15, 313.15, and 323.15) K   |
| dvisc | 0.0007125 | Paxs | 318.15 | Densities, Viscosities, and Sound Speeds of Some Acetate Salts in Binary Mixtures of Tetrahydrofuran and Methanol at (303.15, 313.15, and 323.15) K   |
| dvisc | 0.0007103 | Paxs | 308.15 | Electrical Conductances of Tetrabutylammonium Bromide, Sodium Tetraphenylborate, and Sodium Bromide in N,N-Dimethylformamide (1) + Water (2) Mixtures at (308.15, 313.15, 318.15, and 323.15) K |

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|-------|-----------|------|--------|---|
| dvisc | 0.0006691 | Paxs | 313.15 | Electrical Conductances of Tetrabutylammonium Bromide, Sodium Tetraphenylborate, and Sodium Bromide in N,N-Dimethylformamide (1) + Water (2) Mixtures at (308.15, 313.15, 318.15, and 323.15) K |
| dvisc | 0.0006402 | Paxs | 318.15 | Electrical Conductances of Tetrabutylammonium Bromide, Sodium Tetraphenylborate, and Sodium Bromide in N,N-Dimethylformamide (1) + Water (2) Mixtures at (308.15, 313.15, 318.15, and 323.15) K |
| dvisc | 0.0005999 | Paxs | 323.15 | Electrical Conductances of Tetrabutylammonium Bromide, Sodium Tetraphenylborate, and Sodium Bromide in N,N-Dimethylformamide (1) + Water (2) Mixtures at (308.15, 313.15, 318.15, and 323.15) K |
| dvisc | 0.0010158 | Paxs | 283.15 | Conductometric Studies of 1-Ethyl-3-methylimidazolium Tetrafluoroborate and 1-Butyl-3-methylimidazolium Tetrafluoroborate in N,N-Dimethylformamide at Temperatures from (283.15 to 318.15) K    |

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|-------|-----------|------|--------|---|
| dvisc | 0.0009545 | Paxs | 288.15 | Conductometric<br>Studies of<br>1-Ethyl-3-methylimidazolium<br>Tetrafluoroborate<br>and<br>1-Butyl-3-methylimidazolium<br>Tetrafluoroborate<br>in<br>N,N-Dimethylformamide<br>at Temperatures<br>from (283.15 to<br>318.15) K |
| dvisc | 0.0008985 | Paxs | 293.15 | Conductometric<br>Studies of<br>1-Ethyl-3-methylimidazolium<br>Tetrafluoroborate<br>and<br>1-Butyl-3-methylimidazolium<br>Tetrafluoroborate<br>in<br>N,N-Dimethylformamide<br>at Temperatures<br>from (283.15 to<br>318.15) K |
| dvisc | 0.0008455 | Paxs | 298.15 | Conductometric<br>Studies of<br>1-Ethyl-3-methylimidazolium<br>Tetrafluoroborate<br>and<br>1-Butyl-3-methylimidazolium<br>Tetrafluoroborate<br>in<br>N,N-Dimethylformamide<br>at Temperatures<br>from (283.15 to<br>318.15) K |
| dvisc | 0.0007990 | Paxs | 303.15 | Conductometric<br>Studies of<br>1-Ethyl-3-methylimidazolium<br>Tetrafluoroborate<br>and<br>1-Butyl-3-methylimidazolium<br>Tetrafluoroborate<br>in<br>N,N-Dimethylformamide<br>at Temperatures<br>from (283.15 to<br>318.15) K |

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|-------|-----------|------|--------|---|
| dvisc | 0.0007553 | Paxs | 308.15 | Conductometric<br>Studies of<br>1-Ethyl-3-methylimidazolium<br>Tetrafluoroborate<br>and<br>1-Butyl-3-methylimidazolium<br>Tetrafluoroborate<br>in<br>N,N-Dimethylformamide<br>at Temperatures<br>from (283.15 to<br>318.15) K |
| dvisc | 0.0007172 | Paxs | 313.15 | Conductometric<br>Studies of<br>1-Ethyl-3-methylimidazolium<br>Tetrafluoroborate<br>and<br>1-Butyl-3-methylimidazolium<br>Tetrafluoroborate<br>in<br>N,N-Dimethylformamide<br>at Temperatures<br>from (283.15 to<br>318.15) K |
| dvisc | 0.0006826 | Paxs | 318.15 | Conductometric<br>Studies of<br>1-Ethyl-3-methylimidazolium<br>Tetrafluoroborate<br>and<br>1-Butyl-3-methylimidazolium<br>Tetrafluoroborate<br>in<br>N,N-Dimethylformamide<br>at Temperatures<br>from (283.15 to<br>318.15) K |
| dvisc | 0.0008542 | Paxs | 293.15 | Viscosity of Urea<br>in the Mixture of<br>N,N-Dimethylformamide<br>and Water  |
| dvisc | 0.0005999 | Paxs | 323.15 | Electrical<br>Conductances of<br>Tetrabutylammonium<br>Bromide, Sodium<br>Tetraphenylborate,<br>and Sodium<br>Bromide in<br>N,N-Dimethylformamide<br>at (308.15,<br>313.15, 318.15,<br>and 323.15) K                          |

|       |           |      |        |  |
|-------|-----------|------|--------|--|
| dvisc | 0.0006402 | Paxs | 318.15 | Electrical Conductances of Tetrabutylammonium Bromide, Sodium Tetraphenylborate, and Sodium Bromide in N,N-Dimethylformamide at (308.15, 313.15, 318.15, and 323.15) K |
| dvisc | 0.0006691 | Paxs | 313.15 | Electrical Conductances of Tetrabutylammonium Bromide, Sodium Tetraphenylborate, and Sodium Bromide in N,N-Dimethylformamide at (308.15, 313.15, 318.15, and 323.15) K |
| dvisc | 0.0007103 | Paxs | 308.15 | Electrical Conductances of Tetrabutylammonium Bromide, Sodium Tetraphenylborate, and Sodium Bromide in N,N-Dimethylformamide at (308.15, 313.15, 318.15, and 323.15) K |
| dvisc | 0.0004590 | Paxs | 353.15 | Densities and Viscosities of N,N-Dimethylformamide + N-Methyl-2-pyrrolidinone and + Dimethyl Sulfoxide in the Temperature Range (303.15 to 353.15) K                   |
| dvisc | 0.0005010 | Paxs | 343.15 | Densities and Viscosities of N,N-Dimethylformamide + N-Methyl-2-pyrrolidinone and + Dimethyl Sulfoxide in the Temperature Range (303.15 to 353.15) K                   |
| dvisc | 0.0005520 | Paxs | 333.15 | Densities and Viscosities of N,N-Dimethylformamide + N-Methyl-2-pyrrolidinone and + Dimethyl Sulfoxide in the Temperature Range (303.15 to 353.15) K                   |

|       |           |      |        |  |
|-------|-----------|------|--------|--|
| dvisc | 0.0006080 | Paxs | 323.15 | Densities and Viscosities of N,N-Dimethylformamide + N-Methyl-2-pyrrolidinone and + Dimethyl Sulfoxide in the Temperature Range (303.15 to 353.15) K |
| dvisc | 0.0006780 | Paxs | 313.15 | Densities and Viscosities of N,N-Dimethylformamide + N-Methyl-2-pyrrolidinone and + Dimethyl Sulfoxide in the Temperature Range (303.15 to 353.15) K |
| dvisc | 0.0007590 | Paxs | 303.15 | Densities and Viscosities of N,N-Dimethylformamide + N-Methyl-2-pyrrolidinone and + Dimethyl Sulfoxide in the Temperature Range (303.15 to 353.15) K |
| dvisc | 0.0004600 | Paxs | 353.15 | Densities and Viscosities of N,N-Dimethylformamide + Formic Acid, and + Acetic Acid in the Temperature Range from (303.15 to 353.15) K               |
| dvisc | 0.0005020 | Paxs | 343.15 | Densities and Viscosities of N,N-Dimethylformamide + Formic Acid, and + Acetic Acid in the Temperature Range from (303.15 to 353.15) K               |
| dvisc | 0.0005520 | Paxs | 333.15 | Densities and Viscosities of N,N-Dimethylformamide + Formic Acid, and + Acetic Acid in the Temperature Range from (303.15 to 353.15) K               |

|       |           |      |        |  |
|-------|-----------|------|--------|--|
| dvisc | 0.0006080 | Paxs | 323.15 | Densities and Viscosities of N,N-Dimethylformamide + Formic Acid, and + Acetic Acid in the Temperature Range from (303.15 to 353.15) K |
| dvisc | 0.0006750 | Paxs | 313.15 | Densities and Viscosities of N,N-Dimethylformamide + Formic Acid, and + Acetic Acid in the Temperature Range from (303.15 to 353.15) K |
| dvisc | 0.0007600 | Paxs | 303.15 | Densities and Viscosities of N,N-Dimethylformamide + Formic Acid, and + Acetic Acid in the Temperature Range from (303.15 to 353.15) K |
| dvisc | 0.0004700 | Paxs | 353.15 | Densities and Viscosities of (N,N-Dimethylformamide + Water) at Atmospheric Pressure from (283.15 to 353.15) K                         |
| dvisc | 0.0005100 | Paxs | 343.15 | Densities and Viscosities of (N,N-Dimethylformamide + Water) at Atmospheric Pressure from (283.15 to 353.15) K                         |
| dvisc | 0.0005590 | Paxs | 333.15 | Densities and Viscosities of (N,N-Dimethylformamide + Water) at Atmospheric Pressure from (283.15 to 353.15) K                         |
| dvisc | 0.0006170 | Paxs | 323.15 | Densities and Viscosities of (N,N-Dimethylformamide + Water) at Atmospheric Pressure from (283.15 to 353.15) K                         |

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| dvisc | 0.0006840 | Paxs | 313.15 | Densities and Viscosities of (N,N-Dimethylformamide + Water) at Atmospheric Pressure from (283.15 to 353.15) K  |
| dvisc | 0.0007660 | Paxs | 303.15 | Densities and Viscosities of (N,N-Dimethylformamide + Water) at Atmospheric Pressure from (283.15 to 353.15) K  |
| dvisc | 0.0008640 | Paxs | 293.15 | Densities and Viscosities of (N,N-Dimethylformamide + Water) at Atmospheric Pressure from (283.15 to 353.15) K  |
| dvisc | 0.0005490 | Paxs | 333.15 | Volumetric Properties and Viscosities of Binary Mixtures of N,N-Dimethylformamide with Methanol and Ethanol in the Temperature Range (293.15 to 333.15) K |
| dvisc | 0.0006080 | Paxs | 323.15 | Volumetric Properties and Viscosities of Binary Mixtures of N,N-Dimethylformamide with Methanol and Ethanol in the Temperature Range (293.15 to 333.15) K |
| dvisc | 0.0006750 | Paxs | 313.15 | Volumetric Properties and Viscosities of Binary Mixtures of N,N-Dimethylformamide with Methanol and Ethanol in the Temperature Range (293.15 to 333.15) K |



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| dvisc | 0.0007600 | Paxs   | 303.15 | Volumetric Properties and Viscosities of Binary Mixtures of N,N-Dimethylformamide with Methanol and Ethanol in the Temperature Range (293.15 to 333.15) K |
| dvisc | 0.0008080 | Paxs   | 298.15 | Volumetric Properties and Viscosities of Binary Mixtures of N,N-Dimethylformamide with Methanol and Ethanol in the Temperature Range (293.15 to 333.15) K |
| dvisc | 0.0008630 | Paxs   | 293.15 | Volumetric Properties and Viscosities of Binary Mixtures of N,N-Dimethylformamide with Methanol and Ethanol in the Temperature Range (293.15 to 333.15) K |
| dvisc | 0.0007116 | Paxs   | 308.15 | Viscosity of Urea in the Mixture of N,N-Dimethylformamide and Water   |
| dvisc | 0.0007547 | Paxs   | 303.15 | Viscosity of Urea in the Mixture of N,N-Dimethylformamide and Water   |
| dvisc | 0.0004950 | Paxs   | 338.15 | Densities, Viscosities, and Conductivities of Phosphonic Acid Solutions in N,N-Dimethylformamide and Water  |
| econd | 0.00      | S/m    | 303.15 | Micellar Properties and Related Thermodynamic Parameters of the 14-6-14, 2Br-Gemini Surfactant in Water + Organic Solvent Mixed Media                     |
| hfust | 8.95      | kJ/mol | 212.90 | NIST Webbook  |
| hfust | 8.95      | kJ/mol | 212.90 | NIST Webbook  |
| hfust | 8.95      | kJ/mol | 212.85 | NIST Webbook  |

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|-------|--------|--------|--------|--|
| hfust | 8.95   | kJ/mol | 212.90 | NIST Webbook   |
| hvapt | 46.70  | kJ/mol | 333.00 | NIST Webbook   |
| hvapt | 42.50  | kJ/mol | 370.50 | NIST Webbook   |
| hvapt | 49.20  | kJ/mol | 363.50 | NIST Webbook   |
| hvapt | 43.60  | kJ/mol | 381.50 | NIST Webbook   |
| hvapt | 41.80  | kJ/mol | 401.50 | NIST Webbook   |
| hvapt | 43.10  | kJ/mol | 385.50 | NIST Webbook   |
| hvapt | 56.70  | kJ/mol | 378.00 | NIST Webbook   |
| pvap  | 1.74   | kPa    | 317.15 | Experimental Vapor Pressure Data and a Vapor Pressure Equation for N,N-Dimethylformamide   |
| pvap  | 0.85   | kPa    | 305.20 | Vapour pressures and enthalpies of vaporisation of alkyl formamides  |
| pvap  | 101.30 | kPa    | 426.05 | Separation of azeotropic mixture (2, 2, 3, 3-Tetrafluoro-1-propanol + water) by extractive distillation: Entrainers selection and vapour-liquid equilibrium measurements |
| pvap  | 17.37  | kPa    | 367.15 | Experimental Vapor Pressure Data and a Vapor Pressure Equation for N,N-Dimethylformamide   |
| pvap  | 96.15  | kPa    | 424.60 | Vapor Liquid Equilibrium Data for Binary Mixtures of Acetic Acid + Anisole, Acetone + Anisole, and Isopropanol + Anisole at Pressure 96.15 kPa                           |
| pvap  | 0.92   | kPa    | 307.15 | Experimental Vapor Pressure Data and a Vapor Pressure Equation for N,N-Dimethylformamide   |
| pvap  | 1.06   | kPa    | 309.15 | Experimental Vapor Pressure Data and a Vapor Pressure Equation for N,N-Dimethylformamide   |

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| pvap | 1.20  | kPa | 311.15 | Experimental Vapor Pressure Data and a Vapor Pressure Equation for N,N-Dimethylformamide |
| pvap | 1.36  | kPa | 313.15 | Experimental Vapor Pressure Data and a Vapor Pressure Equation for N,N-Dimethylformamide |
| pvap | 1.53  | kPa | 315.15 | Experimental Vapor Pressure Data and a Vapor Pressure Equation for N,N-Dimethylformamide |
| pvap | 0.69  | kPa | 302.20 | Vapour pressures and enthalpies of vaporisation of alkyl formamides                      |
| pvap | 1.95  | kPa | 319.15 | Experimental Vapor Pressure Data and a Vapor Pressure Equation for N,N-Dimethylformamide |
| pvap | 2.18  | kPa | 321.15 | Experimental Vapor Pressure Data and a Vapor Pressure Equation for N,N-Dimethylformamide |
| pvap | 2.43  | kPa | 323.15 | Experimental Vapor Pressure Data and a Vapor Pressure Equation for N,N-Dimethylformamide |
| pvap | 2.70  | kPa | 325.15 | Experimental Vapor Pressure Data and a Vapor Pressure Equation for N,N-Dimethylformamide |
| pvap | 16.07 | kPa | 365.15 | Experimental Vapor Pressure Data and a Vapor Pressure Equation for N,N-Dimethylformamide |
| pvap | 14.85 | kPa | 363.15 | Experimental Vapor Pressure Data and a Vapor Pressure Equation for N,N-Dimethylformamide |

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| pvap | 0.59   | kPa | 299.50 | Vapour pressures and enthalpies of vaporisation of alkyl formamides   |
| pvap | 12.66  | kPa | 359.15 | Experimental Vapor Pressure Data and a Vapor Pressure Equation for N,N-Dimethylformamide  |
| pvap | 11.67  | kPa | 357.15 | Experimental Vapor Pressure Data and a Vapor Pressure Equation for N,N-Dimethylformamide  |
| pvap | 101.00 | kPa | 425.15 | Vapor Liquid Equilibrium Data for Methanol + tert-Butylamine + N,N-Dimethylformamide and Constituent Binary Systems at Atmospheric Pressure |
| pvap | 10.76  | kPa | 355.15 | Experimental Vapor Pressure Data and a Vapor Pressure Equation for N,N-Dimethylformamide  |
| pvap | 9.90   | kPa | 353.15 | Experimental Vapor Pressure Data and a Vapor Pressure Equation for N,N-Dimethylformamide  |
| pvap | 9.10   | kPa | 351.15 | Experimental Vapor Pressure Data and a Vapor Pressure Equation for N,N-Dimethylformamide  |
| pvap | 13.72  | kPa | 361.15 | Experimental Vapor Pressure Data and a Vapor Pressure Equation for N,N-Dimethylformamide  |
| pvap | 0.48   | kPa | 296.30 | Vapour pressures and enthalpies of vaporisation of alkyl formamides   |
| pvap | 0.40   | kPa | 293.30 | Vapour pressures and enthalpies of vaporisation of alkyl formamides   |

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| pvap | 3.00  | kPa | 327.15 | Experimental Vapor Pressure Data and a Vapor Pressure Equation for N,N-Dimethylformamide |
| pvap | 3.32  | kPa | 329.15 | Experimental Vapor Pressure Data and a Vapor Pressure Equation for N,N-Dimethylformamide |
| pvap | 0.31  | kPa | 289.30 | Vapour pressures and enthalpies of vaporisation of alkyl formamides                      |
| pvap | 18.81 | kPa | 369.15 | Experimental Vapor Pressure Data and a Vapor Pressure Equation for N,N-Dimethylformamide |
| pvap | 5.88  | kPa | 341.15 | Experimental Vapor Pressure Data and a Vapor Pressure Equation for N,N-Dimethylformamide |
| pvap | 0.24  | kPa | 285.30 | Vapour pressures and enthalpies of vaporisation of alkyl formamides                      |
| pvap | 3.67  | kPa | 331.15 | Experimental Vapor Pressure Data and a Vapor Pressure Equation for N,N-Dimethylformamide |
| pvap | 4.04  | kPa | 333.15 | Experimental Vapor Pressure Data and a Vapor Pressure Equation for N,N-Dimethylformamide |
| pvap | 4.45  | kPa | 335.15 | Experimental Vapor Pressure Data and a Vapor Pressure Equation for N,N-Dimethylformamide |
| pvap | 4.89  | kPa | 337.15 | Experimental Vapor Pressure Data and a Vapor Pressure Equation for N,N-Dimethylformamide |

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| pvap | 5.37 | kPa | 339.15 | Experimental Vapor Pressure Data and a Vapor Pressure Equation for N,N-Dimethylformamide |
| pvap | 0.20 | kPa | 283.10 | Vapour pressures and enthalpies of vaporisation of alkyl formamides                      |
| pvap | 6.43 | kPa | 343.15 | Experimental Vapor Pressure Data and a Vapor Pressure Equation for N,N-Dimethylformamide |
| pvap | 7.03 | kPa | 345.15 | Experimental Vapor Pressure Data and a Vapor Pressure Equation for N,N-Dimethylformamide |
| pvap | 7.67 | kPa | 347.15 | Experimental Vapor Pressure Data and a Vapor Pressure Equation for N,N-Dimethylformamide |
| pvap | 8.36 | kPa | 349.15 | Experimental Vapor Pressure Data and a Vapor Pressure Equation for N,N-Dimethylformamide |
| pvap | 0.18 | kPa | 281.20 | Vapour pressures and enthalpies of vaporisation of alkyl formamides                      |
| pvap | 0.15 | kPa | 279.30 | Vapour pressures and enthalpies of vaporisation of alkyl formamides                      |
| pvap | 0.36 | kPa | 291.30 | Vapour pressures and enthalpies of vaporisation of alkyl formamides                      |
| pvap | 0.12 | kPa | 276.30 | Vapour pressures and enthalpies of vaporisation of alkyl formamides                      |

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| pvap | 96.15   | kPa | 424.53 | Vapor Liquid<br>Equilibrium Data<br>for Binary<br>Mixtures of<br>Acetic Acid +<br>Anisole, Acetone<br>+ Anisole, and<br>Isopropanol +<br>Anisole at<br>Pressure 96.15<br>kPa   |
| rfi  | 1.43070 |     | 298.15 | Densities and<br>volumetric<br>properties of<br>binary mixtures<br>of xylene with<br>N,N-dimethylformamide<br>at different<br>temperatures   |
| rfi  | 1.42880 |     | 298.15 | Density,<br>Viscosity,<br>Refractive Index,<br>and Speed of<br>Sound for Binary<br>Mixtures of<br>Anisole with<br>2-Chloroethanol,<br>1,4-Dioxane,<br>Tetrachloroethylene,<br>Tetrachloroethane,<br>DMF, DMSO,<br>and Diethyl<br>Oxalate at<br>(298.15, 303.15,<br>and 308.15) K |
| rfi  | 1.42640 |     | 303.15 | Density,<br>Viscosity,<br>Refractive Index,<br>and Speed of<br>Sound for Binary<br>Mixtures of<br>Anisole with<br>2-Chloroethanol,<br>1,4-Dioxane,<br>Tetrachloroethylene,<br>Tetrachloroethane,<br>DMF, DMSO,<br>and Diethyl<br>Oxalate at<br>(298.15, 303.15,<br>and 308.15) K |

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| rfi | 1.43030 | 293.15 | Solid-Liquid Equilibrium Measurements for Posaconazole and Voriconazole in Several Solvents between T = 278.2 and 323.2 K Using Differential Thermal Analysis/Thermal Gravimetric Analysis |
| rfi | 1.43040 | 293.15 | Experimental solubility for betulin and estrone in various solvents within the temperature range T = (293.2 to 328.2) K  |
| rfi | 1.42900 | 293.15 | Effect of temperature and chain length on the viscosity and surface tension of binary systems of N,N-dimethylformamide with 1-octanol, 1-nonanol and 1-decanol                             |
| rfi | 1.42900 | 293.10 | Viscosity and surface tension of binary systems of N,N-dimethylformamide with alkan-1-ols at different temperatures  |
| rfi | 1.42100 | 313.15 | Steric parameters and excess properties of hydroxamic acids  |
| rfi | 1.42300 | 308.15 | Steric parameters and excess properties of hydroxamic acids  |
| rfi | 1.42530 | 303.15 | Steric parameters and excess properties of hydroxamic acids  |
| rfi | 1.42700 | 298.15 | Steric parameters and excess properties of hydroxamic acids  |



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| rfi | 1.42220 | 313.15 | Comparative study of physico-chemical properties of binary mixtures of N,N-dimethylformamide with 1-alkanols at different temperatures     |
| rfi | 1.42520 | 308.15 | Comparative study of physico-chemical properties of binary mixtures of N,N-dimethylformamide with 1-alkanols at different temperatures     |
| rfi | 1.42710 | 303.15 | Comparative study of physico-chemical properties of binary mixtures of N,N-dimethylformamide with 1-alkanols at different temperatures     |
| rfi | 1.42800 | 298.15 | Comparative study of physico-chemical properties of binary mixtures of N,N-dimethylformamide with 1-alkanols at different temperatures     |
| rfi | 1.43070 | 298.15 | Effects of the presence of ethylacetate or benzene on the densities and volumetric properties of mixture (styrene + N,N-dimethylformamide) |
| rfi | 1.43070 | 298.15 | Densities and excess volumes of binary mixtures of N,N-dimethylformamide with aromatic hydrocarbon at different temperature                |

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| rfi | 1.43090 | 293.15 | Solubility Data for Roflumilast and Maraviroc in Various Solvents between T = (278.2-323.2) K  |
| rfi | 1.42650 | 298.15 | Phase Equilibria Involved in Extractive Distillation of Dipropyl Ether + 1-Propyl Alcohol Using N,N-Dimethylformamide as Entrainer   |
| rfi | 1.43100 | 288.15 | Partial Molar Volumes of N,N'-1,2-Ethyl-bis(salicyladimine) Schiff Base (Salen) in Organic Solvents at T = (283.15 to 318.15) K  |
| rfi | 1.42760 | 298.15 | Partial Molar Volumes of N,N'-1,2-Ethyl-bis(salicyladimine) Schiff Base (Salen) in Organic Solvents at T = (283.15 to 318.15) K  |
| rfi | 1.42300 | 308.15 | Partial Molar Volumes of N,N'-1,2-Ethyl-bis(salicyladimine) Schiff Base (Salen) in Organic Solvents at T = (283.15 to 318.15) K  |
| rfi | 1.41800 | 318.15 | Partial Molar Volumes of N,N'-1,2-Ethyl-bis(salicyladimine) Schiff Base (Salen) in Organic Solvents at T = (283.15 to 318.15) K  |
| rfi | 1.42820 | 298.15 | Vapor-Liquid Equilibrium and Excess Gibbs Energies of Hexane + N,N-Dimethyl Formamide, 2-Methylpropan-2-ol + 2-Aminophenol, N,N-Dimethyl Formamide, and 2-Propanol + Diisopropyl Amine at 94.4 kPa |

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| rfi | 1.43070 | 293.10 | Phase Equilibria of Binary Systems Comprising Formic Acid, N,N-Dimethylformamide, 1-Chloro-2-ethylhexane, and 2-Ethyl-1-hexanol |
| rfi | 1.43070 | 293.10 | Vapor-Liquid Equilibria of Binary Systems Comprising 1-Chloro-2-ethylhexane and 2-Ethyl-1-hexanol                               |
| rfi | 1.42820 | 298.15 | Bubble Temperature Measurements on Binary Mixtures Formed by Cyclohexane at 94.7 kPa  |
| rfi | 1.42190 | 298.15 | Activity coefficients and excess Gibbs energy of binary mixtures of N,N-dimethyl formamide with selected compounds at 95.5 kPa  |
| rfi | 1.43050 | 293.15 | Excess Gibbs' energies of the binary mixtures formed by N,N-dimethylformamide with xylenes and cresols at 95.1 kPa              |
| rfi | 1.42820 | 298.15 | (Vapor + liquid) equilibrium of binary mixtures formed by N,N-dimethyl formamide with some compounds at 95.1 kPa                |
| rfi | 1.43050 | 293.15 | Excess Gibbs energies of binary mixtures formed by nitrobenzene with selected compounds at 94.95 kPa                            |

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| rfi  | 1.42820 |       | 298.15 | Excess Gibbs energies of selected binary mixtures formed by N,N-dimethylformamide at 95.5 kPa   |
| rfi  | 1.42810 |       | 293.10 | Phase equilibria for the extraction of sec-butylbenzene from dodecane with N,N-dimethylformamide  |
| rfi  | 1.42650 |       | 298.15 | Phase equilibria in the systems isobutyl alcohol +N,N-dimethylformamide, isobutyl acetate +N,N-dimethylformamide and isobutyl alcohol + isobutyl acetate +N,N-dimethylformamide at 101.3 kPa  |
| rfi  | 1.42380 |       | 308.15 | Density, Viscosity, Refractive Index, and Speed of Sound for Binary Mixtures of Anisole with 2-Chloroethanol, 1,4-Dioxane, Tetrachloroethylene, Tetrachloroethane, DMF, DMSO, and Diethyl Oxalate at (298.15, 303.15, and 308.15) K |
| rhoI | 943.77  | kg/m3 | 298.15 | Density, Speed of Sound, and Viscosity of Diethylene Glycol Monoethyl Ether + N,N-Dimethylformamide (Ethanol, Water) at T = 288.15-318.15 K   |
| rhoI | 919.86  | kg/m3 | 323.15 | Density, Viscosity, and Electrical Conductivity of Protic Amidium Bis(trifluoromethanesulfonyl)amide Ionic Liquids  |

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|------|--------|-------|--------|---|--|
| rhoI | 910.17 | kg/m3 | 333.15 | Density,<br>Viscosity, and<br>Electrical<br>Conductivity of<br>Protic Amidium<br>Bis(trifluoromethanesulfonyl)amide<br>Ionic Liquids  |  |
| rhoI | 900.41 | kg/m3 | 343.15 | Density,<br>Viscosity, and<br>Electrical<br>Conductivity of<br>Protic Amidium<br>Bis(trifluoromethanesulfonyl)amide<br>Ionic Liquids  |  |
| rhoI | 890.57 | kg/m3 | 353.15 | Density,<br>Viscosity, and<br>Electrical<br>Conductivity of<br>Protic Amidium<br>Bis(trifluoromethanesulfonyl)amide<br>Ionic Liquids  |  |
| rhoI | 880.65 | kg/m3 | 363.15 | Density,<br>Viscosity, and<br>Electrical<br>Conductivity of<br>Protic Amidium<br>Bis(trifluoromethanesulfonyl)amide<br>Ionic Liquids  |  |
| rhoI | 943.89 | kg/m3 | 298.15 | Isobaric Vapor<br>Liquid Equilibria<br>for Two Binary<br>Systems<br>{Propylene<br>Glycol Methyl<br>Ether Acetate +<br>Methanol} and<br>{Propylene<br>Glycol Methyl<br>Ether Acetate +<br>N,N-Dimethylformamide}<br>at p = 30.0, 50.0,<br>and 70.0 kPa |  |
| rhoI | 939.18 | kg/m3 | 303.15 | Influence of<br>Aprotic<br>Cosolvents on<br>the<br>Thermophysical<br>Properties of<br>Imidazolium-Based<br>Ionic Liquid   |  |
| rhoI | 934.41 | kg/m3 | 308.15 | Influence of<br>Aprotic<br>Cosolvents on<br>the<br>Thermophysical<br>Properties of<br>Imidazolium-Based<br>Ionic Liquid   |  |

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| rho1 | 929.64 | kg/m3 | 313.15 | Influence of Aprotic Cosolvents on the Thermophysical Properties of Imidazolium-Based Ionic Liquid |
| rho1 | 924.86 | kg/m3 | 318.15 | Influence of Aprotic Cosolvents on the Thermophysical Properties of Imidazolium-Based Ionic Liquid |
| rho1 | 920.07 | kg/m3 | 323.15 | Influence of Aprotic Cosolvents on the Thermophysical Properties of Imidazolium-Based Ionic Liquid |
| rho1 | 915.26 | kg/m3 | 328.15 | Influence of Aprotic Cosolvents on the Thermophysical Properties of Imidazolium-Based Ionic Liquid |
| rho1 | 910.44 | kg/m3 | 333.15 | Influence of Aprotic Cosolvents on the Thermophysical Properties of Imidazolium-Based Ionic Liquid |
| rho1 | 905.60 | kg/m3 | 338.15 | Influence of Aprotic Cosolvents on the Thermophysical Properties of Imidazolium-Based Ionic Liquid |
| rho1 | 900.74 | kg/m3 | 343.15 | Influence of Aprotic Cosolvents on the Thermophysical Properties of Imidazolium-Based Ionic Liquid |

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| rhoI | 895.87 | kg/m3 | 348.15 | Influence of Aprotic Cosolvents on the Thermophysical Properties of Imidazolium-Based Ionic Liquid  |
| rhoI | 890.98 | kg/m3 | 353.15 | Influence of Aprotic Cosolvents on the Thermophysical Properties of Imidazolium-Based Ionic Liquid  |
| rhoI | 944.69 | kg/m3 | 298.15 | Liquid-Liquid Equilibrium for Ternary Systems, Water + 5-Hydroxymethylfurfural + (1-Butanol, Isobutanol, Methyl Isobutyl Ketone), at 313.15, 323.15, and 333.15 K                         |
| rhoI | 943.93 | kg/m3 | 298.15 | Effect of Solvents and Temperature on Interactions in Binary and Ternary Mixtures of 1-Butyl-3-methylimidazolium Trifluoromethanesulfonate with Acetonitrile or/and N,N-Dimethylformamide |
| rhoI | 939.16 | kg/m3 | 303.15 | Effect of Solvents and Temperature on Interactions in Binary and Ternary Mixtures of 1-Butyl-3-methylimidazolium Trifluoromethanesulfonate with Acetonitrile or/and N,N-Dimethylformamide |
| rhoI | 934.37 | kg/m3 | 308.15 | Effect of Solvents and Temperature on Interactions in Binary and Ternary Mixtures of 1-Butyl-3-methylimidazolium Trifluoromethanesulfonate with Acetonitrile or/and N,N-Dimethylformamide |

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|------|--------|-------|--------|---|
| rhoI | 929.58 | kg/m3 | 313.15 | Effect of Solvents and Temperature on Interactions in Binary and Ternary Mixtures of 1-Butyl-3-methylimidazolium Trifluoromethanesulfonate with Acetonitrile or/and N,N-Dimethylformamide |
| rhoI | 924.76 | kg/m3 | 318.15 | Effect of Solvents and Temperature on Interactions in Binary and Ternary Mixtures of 1-Butyl-3-methylimidazolium Trifluoromethanesulfonate with Acetonitrile or/and N,N-Dimethylformamide |
| rhoI | 919.95 | kg/m3 | 323.15 | Effect of Solvents and Temperature on Interactions in Binary and Ternary Mixtures of 1-Butyl-3-methylimidazolium Trifluoromethanesulfonate with Acetonitrile or/and N,N-Dimethylformamide |
| rhoI | 953.29 | kg/m3 | 288.15 | Density, Speed of Sound, and Viscosity of Diethylene Glycol Monoethyl Ether + N,N-Dimethylformamide (Ethanol, Water) at T = 288.15-318.15 K   |
| rhoI | 929.50 | kg/m3 | 313.15 | Density, Viscosity, and Electrical Conductivity of Protic Amidium Bis(trifluoromethanesulfonyl)amide Ionic Liquids  |
| rhoI | 934.21 | kg/m3 | 308.15 | Density, Speed of Sound, and Viscosity of Diethylene Glycol Monoethyl Ether + N,N-Dimethylformamide (Ethanol, Water) at T = 288.15-318.15 K   |



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| rhoI | 924.59 | kg/m3 | 318.15 | Density, Speed of Sound, and Viscosity of Diethylene Glycol Monoethyl Ether + N,N-Dimethylformamide (Ethanol, Water) at T = 288.15-318.15 K   |
| rhoI | 943.81 | kg/m3 | 298.15 | Liquid-Liquid Equilibrium and Excess Enthalpies in Binary Systems Methylcyclohexane + Methanol and Methylcyclohexane + N,N-Dimethylformamide  |
| rhoI | 943.65 | kg/m3 | 298.15 | Excess Molar Volumes and Kinematic Viscosities for Binary Mixtures of Dipropylene Glycol Monobutyl Ether and Dipropylene Glycol tert-Butyl Ether with 2-Pyrrolidinone, N-Methyl-2-pyrrolidinone, N,N-Dimethylformamide, and N,N-Dimethylacetamide at 298.15 K |
| rhoI | 958.08 | kg/m3 | 283.15 | Thermodynamic Properties of Inorganic Salts in Nonaqueous Solvents. I. Apparent Molar Volumes and Compressibilities of Divalent Transition-Metal Perchlorates in N,N-Dimethylformamide  |
| rhoI | 958.16 | kg/m3 | 283.15 | Thermodynamic Properties of Inorganic Salts in Nonaqueous Solvents. I. Apparent Molar Volumes and Compressibilities of Divalent Transition-Metal Perchlorates in N,N-Dimethylformamide  |

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| rhoI | 958.08 | kg/m3 | 283.15 | Thermodynamic Properties of Inorganic Salts in Nonaqueous Solvents. I. Apparent Molar Volumes and Compressibilities of Divalent Transition-Metal Perchlorates in N,N-Dimethylformamide |
| rhoI | 948.57 | kg/m3 | 293.15 | Thermodynamic Properties of Inorganic Salts in Nonaqueous Solvents. I. Apparent Molar Volumes and Compressibilities of Divalent Transition-Metal Perchlorates in N,N-Dimethylformamide |
| rhoI | 948.65 | kg/m3 | 293.15 | Thermodynamic Properties of Inorganic Salts in Nonaqueous Solvents. I. Apparent Molar Volumes and Compressibilities of Divalent Transition-Metal Perchlorates in N,N-Dimethylformamide |
| rhoI | 943.80 | kg/m3 | 298.15 | Thermodynamic Properties of Inorganic Salts in Nonaqueous Solvents. I. Apparent Molar Volumes and Compressibilities of Divalent Transition-Metal Perchlorates in N,N-Dimethylformamide |
| rhoI | 939.07 | kg/m3 | 303.15 | Density, Viscosity, and Electrical Conductivity of Protic Amidium Bis(trifluoromethanesulfonyl)amide Ionic Liquids   |

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| rhoI | 939.02 | kg/m3 | 303.15 | Thermodynamic Properties of Inorganic Salts in Nonaqueous Solvents. I. Apparent Molar Volumes and Compressibilities of Divalent Transition-Metal Perchlorates in N,N-Dimethylformamide |
| rhoI | 939.11 | kg/m3 | 303.15 | Thermodynamic Properties of Inorganic Salts in Nonaqueous Solvents. I. Apparent Molar Volumes and Compressibilities of Divalent Transition-Metal Perchlorates in N,N-Dimethylformamide |
| rhoI | 929.44 | kg/m3 | 313.15 | Thermodynamic Properties of Inorganic Salts in Nonaqueous Solvents. I. Apparent Molar Volumes and Compressibilities of Divalent Transition-Metal Perchlorates in N,N-Dimethylformamide |
| rhoI | 929.53 | kg/m3 | 313.15 | Thermodynamic Properties of Inorganic Salts in Nonaqueous Solvents. I. Apparent Molar Volumes and Compressibilities of Divalent Transition-Metal Perchlorates in N,N-Dimethylformamide |
| rhoI | 919.81 | kg/m3 | 323.15 | Thermodynamic Properties of Inorganic Salts in Nonaqueous Solvents. I. Apparent Molar Volumes and Compressibilities of Divalent Transition-Metal Perchlorates in N,N-Dimethylformamide |

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| rhoI | 919.90 | kg/m3 | 323.15 | Thermodynamic Properties of Inorganic Salts in Nonaqueous Solvents. I. Apparent Molar Volumes and Compressibilities of Divalent Transition-Metal Perchlorates in N,N-Dimethylformamide |
| rhoI | 910.12 | kg/m3 | 333.15 | Thermodynamic Properties of Inorganic Salts in Nonaqueous Solvents. I. Apparent Molar Volumes and Compressibilities of Divalent Transition-Metal Perchlorates in N,N-Dimethylformamide |
| rhoI | 910.20 | kg/m3 | 333.15 | Thermodynamic Properties of Inorganic Salts in Nonaqueous Solvents. I. Apparent Molar Volumes and Compressibilities of Divalent Transition-Metal Perchlorates in N,N-Dimethylformamide |
| rhoI | 943.81 | kg/m3 | 298.15 | Liquid-Liquid Equilibrium and Excess Enthalpies in the Binary System 2-Methylpentane + N,N-Dimethylformamide   |
| rhoI | 944.20 | kg/m3 | 298.15 | Excess Molar Volumes and Viscosity Deviations for the Ternary System N,N-Dimethylformamide + N-Methylformamide + Water and the Binary Subsystems at 298.15 K                           |

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| rhoI | 943.88 | kg/m3 | 298.15 | Thermodynamics of Mixtures Containing a Strongly Polar Compound. 8. Liquid-Liquid Equilibria for N,N-Dialkylamide + Selected N-Alkanes   |
| rhoI | 943.88 | kg/m3 | 298.15 | Effect of N,N'-Bis(2-pyridylmethylidene)-1,2-diiminoethane (BPIE) Schiff Base on the Thermophysical Properties of Ionic Liquids in N,N-Dimethylformamide Solutions at 298.15 K         |
| rhoI | 948.74 | kg/m3 | 293.15 | Volumetric Properties of Urea in the Mixture of N,N-Dimethylformamide with Water   |
| rhoI | 943.98 | kg/m3 | 298.15 | Volumetric Properties of Urea in the Mixture of N,N-Dimethylformamide with Water   |
| rhoI | 939.20 | kg/m3 | 303.15 | Volumetric Properties of Urea in the Mixture of N,N-Dimethylformamide with Water   |
| rhoI | 934.42 | kg/m3 | 308.15 | Volumetric Properties of Urea in the Mixture of N,N-Dimethylformamide with Water   |
| rhoI | 948.79 | kg/m3 | 293.15 | Volumetric Properties of Binary Mixtures of 1-Butyl-1-methylpyrrolidinium Bis(trifluoromethylsulfonyl)imide with N-Methylformamide and N,N-Dimethylformamide from (293.15 to 323.15) K |

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| rho | 943.87 | kg/m <sup>3</sup> | 298.15 | Volumetric Properties of Binary Mixtures of 1-Butyl-1-methylpyrrolidinium Bis(trifluoromethylsulfonyl)imide with N-Methylformamide and N,N-Dimethylformamide from (293.15 to 323.15) K |
| rho | 938.98 | kg/m <sup>3</sup> | 303.15 | Volumetric Properties of Binary Mixtures of 1-Butyl-1-methylpyrrolidinium Bis(trifluoromethylsulfonyl)imide with N-Methylformamide and N,N-Dimethylformamide from (293.15 to 323.15) K |
| rho | 934.07 | kg/m <sup>3</sup> | 308.15 | Volumetric Properties of Binary Mixtures of 1-Butyl-1-methylpyrrolidinium Bis(trifluoromethylsulfonyl)imide with N-Methylformamide and N,N-Dimethylformamide from (293.15 to 323.15) K |
| rho | 929.11 | kg/m <sup>3</sup> | 313.15 | Volumetric Properties of Binary Mixtures of 1-Butyl-1-methylpyrrolidinium Bis(trifluoromethylsulfonyl)imide with N-Methylformamide and N,N-Dimethylformamide from (293.15 to 323.15) K |
| rho | 924.04 | kg/m <sup>3</sup> | 318.15 | Volumetric Properties of Binary Mixtures of 1-Butyl-1-methylpyrrolidinium Bis(trifluoromethylsulfonyl)imide with N-Methylformamide and N,N-Dimethylformamide from (293.15 to 323.15) K |

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| rhoI | 918.91 | kg/m3 | 323.15 | Volumetric<br>Properties of<br>Binary Mixtures<br>of<br>1-Butyl-1-methylpyrrolidinium<br>Bis(trifluoromethylsulfonyl)imide<br>with<br>N-Methylformamide<br>and<br>N,N-Dimethylformamide<br>from (293.15 to<br>323.15) K   |
| rhoI | 948.72 | kg/m3 | 293.15 | Volumetric<br>Properties of<br>Binary Mixtures<br>of<br>1-Butyl-1-Methylpyrrolidinium<br>Tris(pentafluoroethyl)trifluorophosphate<br>with<br>N-Methylformamide,<br>N-Ethylformamide,<br>N,N-Dimethylformamide,<br>N,N-Dibutylformamide,<br>and<br>N,N-Dimethylacetamide<br>from (293.15 to<br>323.15) K |
| rhoI | 943.86 | kg/m3 | 298.15 | Volumetric<br>Properties of<br>Binary Mixtures<br>of<br>1-Butyl-1-Methylpyrrolidinium<br>Tris(pentafluoroethyl)trifluorophosphate<br>with<br>N-Methylformamide,<br>N-Ethylformamide,<br>N,N-Dimethylformamide,<br>N,N-Dibutylformamide,<br>and<br>N,N-Dimethylacetamide<br>from (293.15 to<br>323.15) K |
| rhoI | 939.00 | kg/m3 | 303.15 | Volumetric<br>Properties of<br>Binary Mixtures<br>of<br>1-Butyl-1-Methylpyrrolidinium<br>Tris(pentafluoroethyl)trifluorophosphate<br>with<br>N-Methylformamide,<br>N-Ethylformamide,<br>N,N-Dimethylformamide,<br>N,N-Dibutylformamide,<br>and<br>N,N-Dimethylacetamide<br>from (293.15 to<br>323.15) K |

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| rhoI | 934.08 | kg/m3 | 308.15 | Volumetric<br>Properties of<br>Binary Mixtures<br>of<br>1-Butyl-1-Methylpyrrolidinium<br>Tris(pentafluoroethyl)trifluorophosphate<br>with<br>N-Methylformamide,<br>N-Ethylformamide,<br>N,N-Dimethylformamide,<br>N,N-Dibutylformamide,<br>and<br>N,N-Dimethylacetamide<br>from (293.15 to<br>323.15) K |
| rhoI | 929.13 | kg/m3 | 313.15 | Volumetric<br>Properties of<br>Binary Mixtures<br>of<br>1-Butyl-1-Methylpyrrolidinium<br>Tris(pentafluoroethyl)trifluorophosphate<br>with<br>N-Methylformamide,<br>N-Ethylformamide,<br>N,N-Dimethylformamide,<br>N,N-Dibutylformamide,<br>and<br>N,N-Dimethylacetamide<br>from (293.15 to<br>323.15) K |
| rhoI | 924.06 | kg/m3 | 318.15 | Volumetric<br>Properties of<br>Binary Mixtures<br>of<br>1-Butyl-1-Methylpyrrolidinium<br>Tris(pentafluoroethyl)trifluorophosphate<br>with<br>N-Methylformamide,<br>N-Ethylformamide,<br>N,N-Dimethylformamide,<br>N,N-Dibutylformamide,<br>and<br>N,N-Dimethylacetamide<br>from (293.15 to<br>323.15) K |
| rhoI | 918.96 | kg/m3 | 323.15 | Volumetric<br>Properties of<br>Binary Mixtures<br>of<br>1-Butyl-1-Methylpyrrolidinium<br>Tris(pentafluoroethyl)trifluorophosphate<br>with<br>N-Methylformamide,<br>N-Ethylformamide,<br>N,N-Dimethylformamide,<br>N,N-Dibutylformamide,<br>and<br>N,N-Dimethylacetamide<br>from (293.15 to<br>323.15) K |



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|------|--------|-------|--------|---|
| rhoI | 953.79 | kg/m3 | 288.15 | Physicochemical Properties for the Binary Systems of Ionic Liquids [Cnmim]Cl + N,N-Dimethylformamide  |
| rhoI | 944.26 | kg/m3 | 298.15 | Physicochemical Properties for the Binary Systems of Ionic Liquids [Cnmim]Cl + N,N-Dimethylformamide  |
| rhoI | 934.69 | kg/m3 | 308.15 | Physicochemical Properties for the Binary Systems of Ionic Liquids [Cnmim]Cl + N,N-Dimethylformamide  |
| rhoI | 925.16 | kg/m3 | 318.15 | Physicochemical Properties for the Binary Systems of Ionic Liquids [Cnmim]Cl + N,N-Dimethylformamide  |
| rhoI | 948.72 | kg/m3 | 293.15 | Volumetric Properties of Binary Mixtures of 1-Butyl-3-Methylimidazolium Tris(pentafluoroethyl)trifluorophosphate with N-Methylformamide, N-Ethylformamide, N,N-Dimethylformamide, N,N-Dibutylformamide, and N,N-Dimethylacetamide from (293.15 to 323.15) K |
| rhoI | 943.86 | kg/m3 | 298.15 | Volumetric Properties of Binary Mixtures of 1-Butyl-3-Methylimidazolium Tris(pentafluoroethyl)trifluorophosphate with N-Methylformamide, N-Ethylformamide, N,N-Dimethylformamide, N,N-Dibutylformamide, and N,N-Dimethylacetamide from (293.15 to 323.15) K |

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|------|--------|-------|--------|---|
| rhoI | 939.00 | kg/m3 | 303.15 | Volumetric<br>Properties of<br>Binary Mixtures<br>of<br>1-Butyl-3-Methylimidazolium<br>Tris(pentafluoroethyl)trifluorophosphate<br>with<br>N-Methylformamide,<br>N-Ethylformamide,<br>N,N-Dimethylformamide,<br>N,N-Dibutylformamide,<br>and<br>N,N-Dimethylacetamide<br>from (293.15 to<br>323.15) K |
| rhoI | 934.08 | kg/m3 | 308.15 | Volumetric<br>Properties of<br>Binary Mixtures<br>of<br>1-Butyl-3-Methylimidazolium<br>Tris(pentafluoroethyl)trifluorophosphate<br>with<br>N-Methylformamide,<br>N-Ethylformamide,<br>N,N-Dimethylformamide,<br>N,N-Dibutylformamide,<br>and<br>N,N-Dimethylacetamide<br>from (293.15 to<br>323.15) K |
| rhoI | 929.13 | kg/m3 | 313.15 | Volumetric<br>Properties of<br>Binary Mixtures<br>of<br>1-Butyl-3-Methylimidazolium<br>Tris(pentafluoroethyl)trifluorophosphate<br>with<br>N-Methylformamide,<br>N-Ethylformamide,<br>N,N-Dimethylformamide,<br>N,N-Dibutylformamide,<br>and<br>N,N-Dimethylacetamide<br>from (293.15 to<br>323.15) K |
| rhoI | 924.06 | kg/m3 | 318.15 | Volumetric<br>Properties of<br>Binary Mixtures<br>of<br>1-Butyl-3-Methylimidazolium<br>Tris(pentafluoroethyl)trifluorophosphate<br>with<br>N-Methylformamide,<br>N-Ethylformamide,<br>N,N-Dimethylformamide,<br>N,N-Dibutylformamide,<br>and<br>N,N-Dimethylacetamide<br>from (293.15 to<br>323.15) K |

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|------|--------|-------|--------|---|
| rhoI | 918.96 | kg/m3 | 323.15 | Volumetric Properties of Binary Mixtures of 1-Butyl-3-Methylimidazolium Tris(pentafluoroethyl)trifluorophosphate with N-Methylformamide, N-Ethylformamide, N,N-Dimethylformamide, N,N-Dibutylformamide, and N,N-Dimethylacetamide from (293.15 to 323.15) K |
| rhoI | 943.82 | kg/m3 | 298.15 | Thermodynamic Properties of Inorganic Salts in Nonaqueous Solvents. III. Apparent Molar Volumes and Compressibilities of Divalent Transition-Metal Chlorides in N,N-Dimethylformamide   |
| rhoI | 943.80 | kg/m3 | 298.15 | Densities and Isothermal Compressibilities at Pressures up to 20 MPa of the Systems N,N-Dimethylformamide or N,N-Dimethylacetamide + r,o-Dichloroalkane   |
| rhoI | 934.20 | kg/m3 | 308.15 | Densities and Isothermal Compressibilities at Pressures up to 20 MPa of the Systems N,N-Dimethylformamide or N,N-Dimethylacetamide + r,o-Dichloroalkane   |
| rhoI | 943.81 | kg/m3 | 298.15 | Surface Tension and Surface Properties of Binary Mixtures of 1,4-Dioxane or N,N-Dimethyl Formamide with n-Alkyl Acetates  |

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|------|--------|-------|--------|---|
| rhoI | 944.21 | kg/m3 | 298.15 | (Liquid + Liquid) Equilibrium for (N,N-Dimethylformamide (DMF) + Hexadecane) at Temperatures between (293.15 and 313.15) K and Ternary Mixtures of (DMF + Hexadecane) with Either Quinoline, or Pyridine, or Pyrrole, or Aniline, or Indole at T = 298.15 K |
| rhoI | 948.63 | kg/m3 | 293.15 | Volumetric Properties of Binary Mixtures of 1-Butyl-3-methylimidazolium Chloride + Water or Hydrophilic Solvents at Different Temperatures  |
| rhoI | 943.86 | kg/m3 | 298.15 | Volumetric Properties of Binary Mixtures of 1-Butyl-3-methylimidazolium Chloride + Water or Hydrophilic Solvents at Different Temperatures  |
| rhoI | 939.08 | kg/m3 | 303.15 | Volumetric Properties of Binary Mixtures of 1-Butyl-3-methylimidazolium Chloride + Water or Hydrophilic Solvents at Different Temperatures  |
| rhoI | 934.29 | kg/m3 | 308.15 | Volumetric Properties of Binary Mixtures of 1-Butyl-3-methylimidazolium Chloride + Water or Hydrophilic Solvents at Different Temperatures  |

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| rhoI | 929.48 | kg/m3 | 313.15 | Volumetric Properties of Binary Mixtures of 1-Butyl-3-methylimidazolium Chloride + Water or Hydrophilic Solvents at Different Temperatures   |  |
| rhoI | 924.68 | kg/m3 | 318.15 | Volumetric Properties of Binary Mixtures of 1-Butyl-3-methylimidazolium Chloride + Water or Hydrophilic Solvents at Different Temperatures   |  |
| rhoI | 943.85 | kg/m3 | 298.15 | Density, Viscosity, and Electrical Conductivity of Protic Amidium Bis(trifluoromethanesulfonyl)amide Ionic Liquids   |  |
| rhoI | 948.61 | kg/m3 | 293.15 | Density, Viscosity, and Electrical Conductivity of Protic Amidium Bis(trifluoromethanesulfonyl)amide Ionic Liquids   |  |
| rhoI | 953.37 | kg/m3 | 288.15 | Density, Viscosity, and Electrical Conductivity of Protic Amidium Bis(trifluoromethanesulfonyl)amide Ionic Liquids   |  |
| rhoI | 943.88 | kg/m3 | 298.15 | Thermodynamic Properties of Inorganic Salts in Nonaqueous Solvents. I. Apparent Molar Volumes and Compressibilities of Divalent Transition-Metal Perchlorates in N,N-Dimethylformamide |  |
| rhoI | 958.10 | kg/m3 | 283.15 | Density, Viscosity, and Electrical Conductivity of Protic Amidium Bis(trifluoromethanesulfonyl)amide Ionic Liquids   |  |

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|------|--------|-------|--------|--|
| rhoI | 962.83 | kg/m3 | 278.15 | Density,<br>Viscosity, and<br>Electrical<br>Conductivity of<br>Protic Amidium<br>Bis(trifluoromethanesulfonyl)amide<br>Ionic Liquids   |
| rhoI | 967.56 | kg/m3 | 273.15 | Density,<br>Viscosity, and<br>Electrical<br>Conductivity of<br>Protic Amidium<br>Bis(trifluoromethanesulfonyl)amide<br>Ionic Liquids   |
| rhoI | 900.57 | kg/m3 | 343.15 | Density and<br>Speed of Sound<br>of Binary<br>Mixtures of Ionic<br>Liquid<br>1-Ethyl-3-methylimidazolium<br>Tetrafluoroborate,<br>N,N-Dimethylformamide,<br>and<br>N,N-Dimethylacetamide<br>at Temperature<br>Range of 293.15<br>343.15 K:<br>Measurement<br>and PC-SAFT<br>Modeling |
| rhoI | 905.45 | kg/m3 | 338.15 | Density and<br>Speed of Sound<br>of Binary<br>Mixtures of Ionic<br>Liquid<br>1-Ethyl-3-methylimidazolium<br>Tetrafluoroborate,<br>N,N-Dimethylformamide,<br>and<br>N,N-Dimethylacetamide<br>at Temperature<br>Range of 293.15<br>343.15 K:<br>Measurement<br>and PC-SAFT<br>Modeling |
| rhoI | 910.31 | kg/m3 | 333.15 | Density and<br>Speed of Sound<br>of Binary<br>Mixtures of Ionic<br>Liquid<br>1-Ethyl-3-methylimidazolium<br>Tetrafluoroborate,<br>N,N-Dimethylformamide,<br>and<br>N,N-Dimethylacetamide<br>at Temperature<br>Range of 293.15<br>343.15 K:<br>Measurement<br>and PC-SAFT<br>Modeling |

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|------|--------|-------|--------|--|
| rhoI | 915.16 | kg/m3 | 328.15 | Density and<br>Speed of Sound<br>of Binary<br>Mixtures of Ionic<br>Liquid<br>1-Ethyl-3-methylimidazolium<br>Tetrafluoroborate,<br>N,N-Dimethylformamide,<br>and<br>N,N-Dimethylacetamide<br>at Temperature<br>Range of 293.15<br>343.15 K:<br>Measurement<br>and PC-SAFT<br>Modeling |
| rhoI | 919.99 | kg/m3 | 323.15 | Density and<br>Speed of Sound<br>of Binary<br>Mixtures of Ionic<br>Liquid<br>1-Ethyl-3-methylimidazolium<br>Tetrafluoroborate,<br>N,N-Dimethylformamide,<br>and<br>N,N-Dimethylacetamide<br>at Temperature<br>Range of 293.15<br>343.15 K:<br>Measurement<br>and PC-SAFT<br>Modeling |
| rhoI | 924.81 | kg/m3 | 318.15 | Density and<br>Speed of Sound<br>of Binary<br>Mixtures of Ionic<br>Liquid<br>1-Ethyl-3-methylimidazolium<br>Tetrafluoroborate,<br>N,N-Dimethylformamide,<br>and<br>N,N-Dimethylacetamide<br>at Temperature<br>Range of 293.15<br>343.15 K:<br>Measurement<br>and PC-SAFT<br>Modeling |
| rhoI | 929.62 | kg/m3 | 313.15 | Density and<br>Speed of Sound<br>of Binary<br>Mixtures of Ionic<br>Liquid<br>1-Ethyl-3-methylimidazolium<br>Tetrafluoroborate,<br>N,N-Dimethylformamide,<br>and<br>N,N-Dimethylacetamide<br>at Temperature<br>Range of 293.15<br>343.15 K:<br>Measurement<br>and PC-SAFT<br>Modeling |

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|------|--------|-------|--------|--|
| rhoI | 934.42 | kg/m3 | 308.15 | Density and<br>Speed of Sound<br>of Binary<br>Mixtures of Ionic<br>Liquid<br>1-Ethyl-3-methylimidazolium<br>Tetrafluoroborate,<br>N,N-Dimethylformamide,<br>and<br>N,N-Dimethylacetamide<br>at Temperature<br>Range of 293.15<br>343.15 K:<br>Measurement<br>and PC-SAFT<br>Modeling |
| rhoI | 939.20 | kg/m3 | 303.15 | Density and<br>Speed of Sound<br>of Binary<br>Mixtures of Ionic<br>Liquid<br>1-Ethyl-3-methylimidazolium<br>Tetrafluoroborate,<br>N,N-Dimethylformamide,<br>and<br>N,N-Dimethylacetamide<br>at Temperature<br>Range of 293.15<br>343.15 K:<br>Measurement<br>and PC-SAFT<br>Modeling |
| rhoI | 943.98 | kg/m3 | 298.15 | Density and<br>Speed of Sound<br>of Binary<br>Mixtures of Ionic<br>Liquid<br>1-Ethyl-3-methylimidazolium<br>Tetrafluoroborate,<br>N,N-Dimethylformamide,<br>and<br>N,N-Dimethylacetamide<br>at Temperature<br>Range of 293.15<br>343.15 K:<br>Measurement<br>and PC-SAFT<br>Modeling |
| rhoI | 948.74 | kg/m3 | 293.15 | Density and<br>Speed of Sound<br>of Binary<br>Mixtures of Ionic<br>Liquid<br>1-Ethyl-3-methylimidazolium<br>Tetrafluoroborate,<br>N,N-Dimethylformamide,<br>and<br>N,N-Dimethylacetamide<br>at Temperature<br>Range of 293.15<br>343.15 K:<br>Measurement<br>and PC-SAFT<br>Modeling |



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|------|--------|-------|--------|---|
| rhoI | 944.08 | kg/m3 | 298.15 | Thermodynamics of amide + amine mixtures. 1. Volumetric, speed of sound, and refractive index data for N,Ndimethylformamide + N-propylpropan-1-amine, + N-butylbutan-1-amine, + butan-1-amine, or + hexan-1-amine systems at several temperatures |
| rhoI | 948.88 | kg/m3 | 293.15 | Thermodynamics of amide + amine mixtures. 1. Volumetric, speed of sound, and refractive index data for N,Ndimethylformamide + N-propylpropan-1-amine, + N-butylbutan-1-amine, + butan-1-amine, or + hexan-1-amine systems at several temperatures |
| rhoI | 910.33 | kg/m3 | 333.15 | Density and Viscosity Measurements for Binary Mixtures of 1-Ethyl-3-methylimidazolium Tetrafluoroborate ([Emim][BF4]) with Dimethylacetamide, Dimethylformamide, and Dimethyl Sulfoxide   |
| rhoI | 920.02 | kg/m3 | 323.15 | Density and Viscosity Measurements for Binary Mixtures of 1-Ethyl-3-methylimidazolium Tetrafluoroborate ([Emim][BF4]) with Dimethylacetamide, Dimethylformamide, and Dimethyl Sulfoxide   |

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| rhoI | 929.64 | kg/m3 | 313.15 | Density and Viscosity Measurements for Binary Mixtures of 1-Ethyl-3-methylimidazolium Tetrafluoroborate ([Emim][BF4]) with Dimethylacetamide, Dimethylformamide, and Dimethyl Sulfoxide        |
| rhoI | 939.23 | kg/m3 | 303.15 | Density and Viscosity Measurements for Binary Mixtures of 1-Ethyl-3-methylimidazolium Tetrafluoroborate ([Emim][BF4]) with Dimethylacetamide, Dimethylformamide, and Dimethyl Sulfoxide        |
| rhoI | 943.70 | kg/m3 | 298.15 | Conductometric, refractometric and FT-IR spectroscopic study of [EMIm]NO3, [EMIm]CH3SO3, and [EMIm]OTs in N,N-dimethyl formamide, N,N-dimethyl acetamide and dimethyl sulphoxide               |
| rhoI | 919.90 | kg/m3 | 323.15 | Excess volumes, Viscosities, and Excess Gibbs Energy of Activation for Viscous Flow, for binary and ternary mixtures 1- propanol + N-N dimethylformamid + chloroform at different temperatures |

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|------|--------|-------|--------|--|
| rhoI | 929.10 | kg/m3 | 313.15 | Excess volumes, Viscosities, and Excess Gibbs Energy of Activation for Viscous Flow, for binary and ternary mixtures 1- propanol + N-N dimethylformamid + chloroform at different temperatures |
| rhoI | 938.70 | kg/m3 | 303.15 | Excess volumes, Viscosities, and Excess Gibbs Energy of Activation for Viscous Flow, for binary and ternary mixtures 1- propanol + N-N dimethylformamid + chloroform at different temperatures |
| rhoI | 948.30 | kg/m3 | 293.15 | Excess volumes, Viscosities, and Excess Gibbs Energy of Activation for Viscous Flow, for binary and ternary mixtures 1- propanol + N-N dimethylformamid + chloroform at different temperatures |
| rhoI | 934.34 | kg/m3 | 308.15 | Thermodynamic evidence for nano-heterogeneity in solutions of the macrocycle C-butylresorcin[4]arene in non-aqueous solvents   |
| rhoI | 939.12 | kg/m3 | 303.15 | Thermodynamic evidence for nano-heterogeneity in solutions of the macrocycle C-butylresorcin[4]arene in non-aqueous solvents   |

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| rhoI | 943.89 | kg/m3 | 298.15 | Thermodynamic evidence for nano-heterogeneity in solutions of the macrocycle C-butylresorcin[4]arene in non-aqueous solvents   |
| rhoI | 948.66 | kg/m3 | 293.15 | Thermodynamic evidence for nano-heterogeneity in solutions of the macrocycle C-butylresorcin[4]arene in non-aqueous solvents   |
| rhoI | 953.41 | kg/m3 | 288.15 | Thermodynamic evidence for nano-heterogeneity in solutions of the macrocycle C-butylresorcin[4]arene in non-aqueous solvents   |
| rhoI | 958.17 | kg/m3 | 283.15 | Thermodynamic evidence for nano-heterogeneity in solutions of the macrocycle C-butylresorcin[4]arene in non-aqueous solvents   |
| rhoI | 962.91 | kg/m3 | 278.15 | Thermodynamic evidence for nano-heterogeneity in solutions of the macrocycle C-butylresorcin[4]arene in non-aqueous solvents   |
| rhoI | 924.73 | kg/m3 | 318.15 | Solvation of ionic liquids based on N-methyl-N-alkylmorpholinium cations in N,N-dimethylformamide and N,N-dimethylacetamide - Volumetric and compressibility studies |
| rhoI | 929.54 | kg/m3 | 313.15 | Solvation of ionic liquids based on N-methyl-N-alkylmorpholinium cations in N,N-dimethylformamide and N,N-dimethylacetamide - Volumetric and compressibility studies |

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|------|--------|-------|--------|--|
| rhoI | 934.34 | kg/m3 | 308.15 | Solvation of ionic liquids based on N-methyl-N-alkylmorpholinium cations in N,N-dimethylformamide and N,N-dimethylacetamide - Volumetric and compressibility studies |
| rhoI | 939.12 | kg/m3 | 303.15 | Solvation of ionic liquids based on N-methyl-N-alkylmorpholinium cations in N,N-dimethylformamide and N,N-dimethylacetamide - Volumetric and compressibility studies |
| rhoI | 943.90 | kg/m3 | 298.15 | Solvation of ionic liquids based on N-methyl-N-alkylmorpholinium cations in N,N-dimethylformamide and N,N-dimethylacetamide - Volumetric and compressibility studies |
| rhoI | 890.98 | kg/m3 | 353.15 | Effect of organic solvents on lowering the viscosity of 1-hexyl-3-methylimidazolium chloride   |
| rhoI | 895.87 | kg/m3 | 348.15 | Effect of organic solvents on lowering the viscosity of 1-hexyl-3-methylimidazolium chloride   |
| rhoI | 900.74 | kg/m3 | 343.15 | Effect of organic solvents on lowering the viscosity of 1-hexyl-3-methylimidazolium chloride   |
| rhoI | 905.60 | kg/m3 | 338.15 | Effect of organic solvents on lowering the viscosity of 1-hexyl-3-methylimidazolium chloride   |

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|------|--------|-------|--------|---|
| rhoI | 910.44 | kg/m3 | 333.15 | Effect of organic solvents on lowering the viscosity of 1-hexyl-3-methylimidazolium chloride  |
| rhoI | 915.26 | kg/m3 | 328.15 | Effect of organic solvents on lowering the viscosity of 1-hexyl-3-methylimidazolium chloride  |
| rhoI | 920.07 | kg/m3 | 323.15 | Effect of organic solvents on lowering the viscosity of 1-hexyl-3-methylimidazolium chloride  |
| rhoI | 924.86 | kg/m3 | 318.15 | Effect of organic solvents on lowering the viscosity of 1-hexyl-3-methylimidazolium chloride  |
| rhoI | 929.64 | kg/m3 | 313.15 | Effect of organic solvents on lowering the viscosity of 1-hexyl-3-methylimidazolium chloride  |
| rhoI | 934.41 | kg/m3 | 308.15 | Effect of organic solvents on lowering the viscosity of 1-hexyl-3-methylimidazolium chloride  |
| rhoI | 939.18 | kg/m3 | 303.15 | Effect of organic solvents on lowering the viscosity of 1-hexyl-3-methylimidazolium chloride  |
| rhoI | 915.14 | kg/m3 | 328.15 | Solvation of ionic liquids based on N-alkyl-N-methylmorpholinium cations in N,N-dimethylformamide and dimethyl sulfoxide. A volumetric and acoustic study |

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|------|--------|-------|--------|---|
| rhoI | 924.80 | kg/m3 | 318.15 | Solvation of ionic liquids based on N-alkyl-N-methylmorpholinium cations in N,N-dimethylformamide and dimethyl sulfoxide. A volumetric and acoustic study                         |
| rhoI | 929.61 | kg/m3 | 313.15 | Solvation of ionic liquids based on N-alkyl-N-methylmorpholinium cations in N,N-dimethylformamide and dimethyl sulfoxide. A volumetric and acoustic study                         |
| rhoI | 934.40 | kg/m3 | 308.15 | Solvation of ionic liquids based on N-alkyl-N-methylmorpholinium cations in N,N-dimethylformamide and dimethyl sulfoxide. A volumetric and acoustic study                         |
| rhoI | 939.19 | kg/m3 | 303.15 | Solvation of ionic liquids based on N-alkyl-N-methylmorpholinium cations in N,N-dimethylformamide and dimethyl sulfoxide. A volumetric and acoustic study                         |
| rhoI | 943.95 | kg/m3 | 298.15 | Solvation of ionic liquids based on N-alkyl-N-methylmorpholinium cations in N,N-dimethylformamide and dimethyl sulfoxide. A volumetric and acoustic study                         |
| rhoI | 935.00 | kg/m3 | 308.00 | Densities, ultrasonic speeds and refractive indices of phenetole with N-methyl-2-pyrrolidone, N,N-dimethylformamide and tetrahydrofuran binary mixtures at different temperatures |

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|------|--------|-------|--------|---|
| rhoI | 939.00 | kg/m3 | 303.00 | Densities, ultrasonic speeds and refractive indices of phenetole with N-methyl-2-pyrrolidone, N,N-dimethylformamide and tetrahydrofuran binary mixtures at different temperatures |
| rhoI | 944.00 | kg/m3 | 298.00 | Densities, ultrasonic speeds and refractive indices of phenetole with N-methyl-2-pyrrolidone, N,N-dimethylformamide and tetrahydrofuran binary mixtures at different temperatures |
| rhoI | 939.28 | kg/m3 | 303.15 | Thermodynamics of amide + ketone mixtures. 1. Volumetric, speed of sound and refractive index data for N,N-dimethylformamide + 2-alkanone systems at several temperatures         |
| rhoI | 944.06 | kg/m3 | 298.15 | Thermodynamics of amide + ketone mixtures. 1. Volumetric, speed of sound and refractive index data for N,N-dimethylformamide + 2-alkanone systems at several temperatures         |
| rhoI | 948.82 | kg/m3 | 293.15 | Thermodynamics of amide + ketone mixtures. 1. Volumetric, speed of sound and refractive index data for N,N-dimethylformamide + 2-alkanone systems at several temperatures         |



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| rhoI | 924.74 | kg/m3 | 318.15 | Solvation of alkaline earth metal ions in N,N-dimethylformamide and N,N-dimethylacetamide - A volumetric and acoustic study  |
| rhoI | 929.55 | kg/m3 | 313.15 | Solvation of alkaline earth metal ions in N,N-dimethylformamide and N,N-dimethylacetamide - A volumetric and acoustic study  |
| rhoI | 934.35 | kg/m3 | 308.15 | Solvation of alkaline earth metal ions in N,N-dimethylformamide and N,N-dimethylacetamide - A volumetric and acoustic study  |
| rhoI | 939.14 | kg/m3 | 303.15 | Solvation of alkaline earth metal ions in N,N-dimethylformamide and N,N-dimethylacetamide - A volumetric and acoustic study  |
| rhoI | 943.91 | kg/m3 | 298.15 | Solvation of alkaline earth metal ions in N,N-dimethylformamide and N,N-dimethylacetamide - A volumetric and acoustic study  |
| rhoI | 935.05 | kg/m3 | 308.15 | Excess molar volumes and excess isentropic compressibilities of binary and ternary mixtures of o-chlorotoluene with cyclic ether and amides or cyclohexane at different temperatures |

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|------|--------|-------|--------|--|
| rhoI | 939.83 | kg/m3 | 303.15 | Excess molar volumes and excess isentropic compressibilities of binary and ternary mixtures of o-chlorotoluene with cyclic ether and amides or cyclohexane at different temperatures |
| rhoI | 944.60 | kg/m3 | 298.15 | Excess molar volumes and excess isentropic compressibilities of binary and ternary mixtures of o-chlorotoluene with cyclic ether and amides or cyclohexane at different temperatures |
| rhoI | 934.39 | kg/m3 | 308.15 | Solution behavior of<br>{(formamide/N-methylformamide/<br>N,N-dimethylformamide)<br>+ CsCl + water}<br>ternary systems<br>at multiple<br>temperatures                                |
| rhoI | 943.89 | kg/m3 | 298.15 | Solution behavior of<br>{(formamide/N-methylformamide/<br>N,N-dimethylformamide)<br>+ CsCl + water}<br>ternary systems<br>at multiple<br>temperatures                                |
| rhoI | 953.53 | kg/m3 | 288.15 | Solution behavior of<br>{(formamide/N-methylformamide/<br>N,N-dimethylformamide)<br>+ CsCl + water}<br>ternary systems<br>at multiple<br>temperatures                                |
| rhoI | 943.97 | kg/m3 | 298.15 | Ionic molar volumes in methanol mixtures with acetonitrile, N,N-dimethylformamide and propylene carbonate at T = 298.15 K  |

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|------|--------|-------|--------|--|
| rhoI | 929.91 | kg/m3 | 313.15 | Solution thermodynamics of iron(III)-N,N'-ethylene-bis(salicylideneiminato)-chloride in binary mixtures of N,N-dimethylformamide and acetonitrile at T = (298.15, 303.15, 308.15 and 313.15) K |
| rhoI | 935.71 | kg/m3 | 308.15 | Solution thermodynamics of iron(III)-N,N'-ethylene-bis(salicylideneiminato)-chloride in binary mixtures of N,N-dimethylformamide and acetonitrile at T = (298.15, 303.15, 308.15 and 313.15) K |
| rhoI | 940.05 | kg/m3 | 303.15 | Solution thermodynamics of iron(III)-N,N'-ethylene-bis(salicylideneiminato)-chloride in binary mixtures of N,N-dimethylformamide and acetonitrile at T = (298.15, 303.15, 308.15 and 313.15) K |
| rhoI | 944.65 | kg/m3 | 298.15 | Solution thermodynamics of iron(III)-N,N'-ethylene-bis(salicylideneiminato)-chloride in binary mixtures of N,N-dimethylformamide and acetonitrile at T = (298.15, 303.15, 308.15 and 313.15) K |
| rhoI | 935.05 | kg/m3 | 308.15 | Topological and thermodynamic investigations of mixtures containing o-chlorotoluene and lower amides   |
| rhoI | 939.83 | kg/m3 | 303.15 | Topological and thermodynamic investigations of mixtures containing o-chlorotoluene and lower amides   |

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|------|--------|-------|--------|---|
| rhoI | 944.60 | kg/m3 | 298.15 | Topological and thermodynamic investigations of mixtures containing o-chlorotoluene and lower amides                                  |
| rhoI | 933.60 | kg/m3 | 313.00 | Thermo-acoustical studies of 1,3,4-oxadiazole as binary mixture at three different temperatures                                       |
| rhoI | 937.20 | kg/m3 | 308.00 | Thermo-acoustical studies of 1,3,4-oxadiazole as binary mixture at three different temperatures                                       |
| rhoI | 941.70 | kg/m3 | 303.00 | Thermo-acoustical studies of 1,3,4-oxadiazole as binary mixture at three different temperatures                                       |
| rhoI | 924.80 | kg/m3 | 318.15 | Solubility for dilute sulfur dioxide in binary mixtures of N,N-dimethylformamide + Ethylene Glycol at T = 308.15 K and p = 122.66 kPa |
| rhoI | 929.50 | kg/m3 | 313.15 | Solubility for dilute sulfur dioxide in binary mixtures of N,N-dimethylformamide + Ethylene Glycol at T = 308.15 K and p = 122.66 kPa |
| rhoI | 934.50 | kg/m3 | 308.15 | Solubility for dilute sulfur dioxide in binary mixtures of N,N-dimethylformamide + Ethylene Glycol at T = 308.15 K and p = 122.66 kPa |

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|------|--------|-------|--------|--|
| rhoI | 939.30 | kg/m3 | 303.15 | Solubility for dilute sulfur dioxide in binary mixtures of N,N-dimethylformamide + Ethylene Glycol at T = 308.15 K and p = 122.66 kPa  |
| rhoI | 944.10 | kg/m3 | 298.15 | Solubility for dilute sulfur dioxide in binary mixtures of N,N-dimethylformamide + Ethylene Glycol at T = 308.15 K and p = 122.66 kPa  |
| rhoI | 926.60 | kg/m3 | 318.15 | Densities, ultrasonic speeds, viscosities and excess properties of binary mixtures of methyl methacrylate with N,N-dimethylformamide and N,N-dimethylacetamide at different temperatures |
| rhoI | 931.10 | kg/m3 | 313.15 | Densities, ultrasonic speeds, viscosities and excess properties of binary mixtures of methyl methacrylate with N,N-dimethylformamide and N,N-dimethylacetamide at different temperatures |
| rhoI | 935.60 | kg/m3 | 308.15 | Densities, ultrasonic speeds, viscosities and excess properties of binary mixtures of methyl methacrylate with N,N-dimethylformamide and N,N-dimethylacetamide at different temperatures |

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|------|--------|-------|--------|---|
| rhoI | 940.10 | kg/m3 | 303.15 | Densities,<br>ultrasonic<br>speeds,<br>viscosities and<br>excess properties<br>of binary<br>mixtures of<br>methyl<br>methacrylate with<br>N,N-dimethylformamide<br>and<br>N,N-dimethylacetamide<br>at different<br>temperatures |
| rhoI | 944.60 | kg/m3 | 298.15 | Densities,<br>ultrasonic<br>speeds,<br>viscosities and<br>excess properties<br>of binary<br>mixtures of<br>methyl<br>methacrylate with<br>N,N-dimethylformamide<br>and<br>N,N-dimethylacetamide<br>at different<br>temperatures |
| rhoI | 949.10 | kg/m3 | 293.15 | Densities,<br>ultrasonic<br>speeds,<br>viscosities and<br>excess properties<br>of binary<br>mixtures of<br>methyl<br>methacrylate with<br>N,N-dimethylformamide<br>and<br>N,N-dimethylacetamide<br>at different<br>temperatures |
| rhoI | 953.61 | kg/m3 | 288.15 | Densities,<br>ultrasonic<br>speeds,<br>viscosities and<br>excess properties<br>of binary<br>mixtures of<br>methyl<br>methacrylate with<br>N,N-dimethylformamide<br>and<br>N,N-dimethylacetamide<br>at different<br>temperatures |

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|------|--------|-------|--------|--|
| rhoI | 944.29 | kg/m3 | 298.15 | Physico-chemical properties of binary mixtures of N,N-dimethylformamide with 1-octanol, 1-nonanol and 1-decanol at different temperatures                                |
| rhoI | 910.11 | kg/m3 | 333.15 | Apparent molar volumes and compressibilities of lanthanum, gadolinium, lutetium and sodium trifluoromethanesulfonates in N,N-dimethylformamide and N,N-dimethylacetamide |
| rhoI | 919.81 | kg/m3 | 323.15 | Apparent molar volumes and compressibilities of lanthanum, gadolinium, lutetium and sodium trifluoromethanesulfonates in N,N-dimethylformamide and N,N-dimethylacetamide |
| rhoI | 929.43 | kg/m3 | 313.15 | Apparent molar volumes and compressibilities of lanthanum, gadolinium, lutetium and sodium trifluoromethanesulfonates in N,N-dimethylformamide and N,N-dimethylacetamide |
| rhoI | 939.01 | kg/m3 | 303.15 | Apparent molar volumes and compressibilities of lanthanum, gadolinium, lutetium and sodium trifluoromethanesulfonates in N,N-dimethylformamide and N,N-dimethylacetamide |

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| rhoI | 943.79 | kg/m3 | 298.15 | Apparent molar volumes and compressibilities of lanthanum, gadolinium, lutetium and sodium trifluoromethanesulfonates in N,N-dimethylformamide and N,N-dimethylacetamide  |
| rhoI | 948.56 | kg/m3 | 293.15 | Apparent molar volumes and compressibilities of lanthanum, gadolinium, lutetium and sodium trifluoromethanesulfonates in N,N-dimethylformamide and N,N-dimethylacetamide  |
| rhoI | 958.07 | kg/m3 | 283.15 | Apparent molar volumes and compressibilities of lanthanum, gadolinium, lutetium and sodium trifluoromethanesulfonates in N,N-dimethylformamide and N,N-dimethylacetamide  |
| rhoI | 944.60 | kg/m3 | 298.15 | Densities and volumetric properties of binary mixtures of N,N-dimethylformamide/N,N-dimethylacetamide with some alkyl acrylates at temperatures from 288.15 K to 318.15 K |
| rhoI | 933.60 | kg/m3 | 313.00 | Thermo-physical properties of 1,3,4-oxadiazole derivatives in pure solvents   |
| rhoI | 937.20 | kg/m3 | 308.00 | Thermo-physical properties of 1,3,4-oxadiazole derivatives in pure solvents   |
| rhoI | 941.70 | kg/m3 | 303.00 | Thermo-physical properties of 1,3,4-oxadiazole derivatives in pure solvents   |



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|------|--------|-------|--------|--|
| rhoI | 944.50 | kg/m3 | 298.15 | Physics and chemistry of an ionic liquid in some industrially important solvent media probed by physicochemical techniques   |
| rhoI | 924.68 | kg/m3 | 318.15 | Apparent molar volumes and expansibilities of H2O and D2O in N,N-dimethylformamide and N,N-dimethylacetamide in the range of T = (278.15 to 318.15) K at p = 0.1 MPa: A comparative analysis |
| rhoI | 934.30 | kg/m3 | 308.15 | Apparent molar volumes and expansibilities of H2O and D2O in N,N-dimethylformamide and N,N-dimethylacetamide in the range of T = (278.15 to 318.15) K at p = 0.1 MPa: A comparative analysis |
| rhoI | 943.87 | kg/m3 | 298.15 | Apparent molar volumes and expansibilities of H2O and D2O in N,N-dimethylformamide and N,N-dimethylacetamide in the range of T = (278.15 to 318.15) K at p = 0.1 MPa: A comparative analysis |
| rhoI | 953.40 | kg/m3 | 288.15 | Apparent molar volumes and expansibilities of H2O and D2O in N,N-dimethylformamide and N,N-dimethylacetamide in the range of T = (278.15 to 318.15) K at p = 0.1 MPa: A comparative analysis |

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|------|--------|-------|--------|--|
| rhoI | 962.90 | kg/m3 | 278.15 | Apparent molar volumes and expansibilities of H2O and D2O in N,N-dimethylformamide and N,N-dimethylacetamide in the range of T = (278.15 to 318.15) K at p = 0.1 MPa: A comparative analysis |
| rhoI | 944.34 | kg/m3 | 298.15 | Excess molar enthalpies of binary systems containing 2-octanone, hexanoic acid, or octanoic acid at T = 298.15 K   |
| rhoI | 943.81 | kg/m3 | 298.15 | Excess molar enthalpies and (vapour + liquid) equilibria for mixtures containing N,N-dialkylamides and alpha,x-dichloroalkanes   |
| rhoI | 943.81 | kg/m3 | 298.15 | (Vapour + liquid) equilibria and excess molar enthalpies for binary mixtures containing N,N-dialkylamides and 1-chloroalkanes  |
| rhoI | 935.55 | kg/m3 | 308.15 | Structural and interactional studies of homologous series of alpha,x-alkanediols in N,N-dimethylformamide  |
| rhoI | 926.70 | kg/m3 | 318.15 | Densities and volumetric properties of (acetonitrile + an amide) binary mixtures at temperatures between 293.15 K and 318.15 K   |
| rhoI | 931.20 | kg/m3 | 313.15 | Densities and volumetric properties of (acetonitrile + an amide) binary mixtures at temperatures between 293.15 K and 318.15 K   |

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|------|--------|-------|--------|---|--|
| rhoI | 935.70 | kg/m3 | 308.15 | Densities and volumetric properties of (acetonitrile + an amide) binary mixtures at temperatures between 293.15 K and 318.15 K              |  |
| rhoI | 940.10 | kg/m3 | 303.15 | Densities and volumetric properties of (acetonitrile + an amide) binary mixtures at temperatures between 293.15 K and 318.15 K              |  |
| rhoI | 944.60 | kg/m3 | 298.15 | Densities and volumetric properties of (acetonitrile + an amide) binary mixtures at temperatures between 293.15 K and 318.15 K              |  |
| rhoI | 949.10 | kg/m3 | 293.15 | Densities and volumetric properties of (acetonitrile + an amide) binary mixtures at temperatures between 293.15 K and 318.15 K              |  |
| rhoI | 935.55 | kg/m3 | 308.15 | A comparative study of thermophysical and spectroscopic properties in mixtures of isomeric butanediol and N,N-dimethylformamide             |  |
| rhoI | 949.17 | kg/m3 | 293.15 | Volumetric properties of binary mixtures of (water + organic solvents) at temperatures between T = 288.15 K and T = 303.15 K at p = 0.1 MPa |  |

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|------|--------|-------|--------|---|
| rhoI | 943.79 | kg/m3 | 298.15 | Limiting partial molar volumes and expansibilities of ammonium perchlorate, tetraalkylammonium perchlorates, and tetrabutylammonium tetraphenylborate in N,N-dimethylformamide          |
| rhoI | 885.68 | kg/m3 | 358.15 | Volumetric properties of (N,N-dimethylformamide + aliphatic diethers) at Temperatures ranging from (298.15 to 358.15) K   |
| rhoI | 900.41 | kg/m3 | 343.15 | Volumetric properties of (N,N-dimethylformamide + aliphatic diethers) at Temperatures ranging from (298.15 to 358.15) K   |
| rhoI | 930.98 | kg/m3 | 313.15 | Volumetric properties of (N,N-dimethylformamide + aliphatic diethers) at Temperatures ranging from (298.15 to 358.15) K   |
| rhoI | 943.85 | kg/m3 | 298.15 | Volumetric properties of (N,N-dimethylformamide + aliphatic diethers) at Temperatures ranging from (298.15 to 358.15) K   |
| rhoI | 943.98 | kg/m3 | 298.15 | PrhoT measurement and PC-SAFT modeling of N,N-dimethyl formamide, N-methyl formamide, N,N-dimethyl acetamide, and ethylenediamine from T = (293.15-423.15) K and pressures up to 35 MPa |

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|------|--------|-------|--------|---|
| rhoI | 924.70 | kg/m3 | 318.15 | Solubility for dilute sulfur dioxide, viscosities, excess properties, and viscous flow thermodynamics of binary system N, N-dimethylformamide + diethylene glycol |
| rhoI | 929.90 | kg/m3 | 313.15 | Solubility for dilute sulfur dioxide, viscosities, excess properties, and viscous flow thermodynamics of binary system N, N-dimethylformamide + diethylene glycol |
| rhoI | 934.20 | kg/m3 | 308.15 | Solubility for dilute sulfur dioxide, viscosities, excess properties, and viscous flow thermodynamics of binary system N, N-dimethylformamide + diethylene glycol |
| rhoI | 939.60 | kg/m3 | 303.15 | Solubility for dilute sulfur dioxide, viscosities, excess properties, and viscous flow thermodynamics of binary system N, N-dimethylformamide + diethylene glycol |
| rhoI | 943.90 | kg/m3 | 298.15 | Solubility for dilute sulfur dioxide, viscosities, excess properties, and viscous flow thermodynamics of binary system N, N-dimethylformamide + diethylene glycol |

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|------|--------|-------|--------|--|--|
| rhoI | 943.83 | kg/m3 | 298.15 | Volumetric investigation of the ternary system ethanol + dimethylformamide + cyclohexane at 298.15 K   |  |
| rhoI | 948.70 | kg/m3 | 293.15 | Bubble point measurements of binary mixtures formed by 1-hexanol with selected nitro-compounds and substituted benzenes at 95.6 kPa                      |  |
| rhoI | 916.40 | kg/m3 | 313.15 | Effect of temperature on ultrasonic velocity and thermodynamic parameters of bisphenol-C-formaldehyde-acrylate resin solutions                           |  |
| rhoI | 919.70 | kg/m3 | 308.15 | Effect of temperature on ultrasonic velocity and thermodynamic parameters of bisphenol-C-formaldehyde-acrylate resin solutions                           |  |
| rhoI | 927.20 | kg/m3 | 303.15 | Effect of temperature on ultrasonic velocity and thermodynamic parameters of bisphenol-C-formaldehyde-acrylate resin solutions                           |  |
| rhoI | 943.81 | kg/m3 | 298.15 | Liquid liquid equilibrium in ternary systems N,N-dimethylformamide + 2-methylpentane + methanol and N,N-dimethylformamide + methylcyclohexane + methanol |  |
| rhoI | 943.81 | kg/m3 | 298.15 | Revision of the volumetric method for measurements of liquid liquid equilibria in binary systems   |  |

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|------|--------|-------|--------|--|
| rhoI | 932.50 | kg/m3 | 313.15 | Interpretation of Association Behavior and Molecular Interactions in Binary Mixtures from Thermoacoustics and Molecular Compression Data |
| rhoI | 935.70 | kg/m3 | 308.15 | Interpretation of Association Behavior and Molecular Interactions in Binary Mixtures from Thermoacoustics and Molecular Compression Data |
| rhoI | 941.90 | kg/m3 | 303.15 | Interpretation of Association Behavior and Molecular Interactions in Binary Mixtures from Thermoacoustics and Molecular Compression Data |
| rhoI | 950.10 | kg/m3 | 298.15 | Interpretation of Association Behavior and Molecular Interactions in Binary Mixtures from Thermoacoustics and Molecular Compression Data |
| rhoI | 955.10 | kg/m3 | 293.15 | Interpretation of Association Behavior and Molecular Interactions in Binary Mixtures from Thermoacoustics and Molecular Compression Data |

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|---------|---------|---------|--------|---|
| rhoI    | 943.78  | kg/m3   | 298.15 | Thermodynamics of amide + amine mixtures. 5. Excess molar enthalpies of N,N-dimethylformamide or N,N-dimethylacetamide + N-propylpropan-1-amine, + N-butylbutan-1-amine, + butan-1-amine, or + hexan-1-amine systems at 298.15 K. Application of the ERAS model |
| rhoI    | 939.36  | kg/m3   | 303.15 | Thermodynamics of amide + amine mixtures. 1. Volumetric, speed of sound, and refractive index data for N,Ndimethylformamide + N-propylpropan-1-amine, + N-butylbutan-1-amine, + butan-1-amine, or + hexan-1-amine systems at several temperatures               |
| rhoI    | 915.07  | kg/m3   | 328.15 | Volumetric properties of (N,N-dimethylformamide + aliphatic diethers) at Temperatures ranging from (298.15 to 358.15) K   |
| sfust   | 41.80   | J/molxK | 212.85 | NIST Webbook  |
| speedsl | 1457.13 | m/s     | 298.15 | Apparent Molar Compressibilities and Volumes of Some 1,1-Electrolytes in N,N-Dimethylacetamide and N,N-Dimethylformamide  |



|         |         |     |        |  |
|---------|---------|-----|--------|--|
| speedsl | 1496.60 | m/s | 288.15 | Volumetric and Isentropic Compressibility Behavior of Ionic Liquid, 1-Propyl-3-Methylimidazolium Bromide in Acetonitrile, Dimethylformamide, and Dimethylsulfoxide at T = (288.15 to 308.15) K |
| speedsl | 1477.27 | m/s | 293.15 | Volumetric and Isentropic Compressibility Behavior of Ionic Liquid, 1-Propyl-3-Methylimidazolium Bromide in Acetonitrile, Dimethylformamide, and Dimethylsulfoxide at T = (288.15 to 308.15) K |
| speedsl | 1457.81 | m/s | 298.15 | Volumetric and Isentropic Compressibility Behavior of Ionic Liquid, 1-Propyl-3-Methylimidazolium Bromide in Acetonitrile, Dimethylformamide, and Dimethylsulfoxide at T = (288.15 to 308.15) K |
| speedsl | 1438.38 | m/s | 303.15 | Volumetric and Isentropic Compressibility Behavior of Ionic Liquid, 1-Propyl-3-Methylimidazolium Bromide in Acetonitrile, Dimethylformamide, and Dimethylsulfoxide at T = (288.15 to 308.15) K |
| speedsl | 1419.04 | m/s | 308.15 | Volumetric and Isentropic Compressibility Behavior of Ionic Liquid, 1-Propyl-3-Methylimidazolium Bromide in Acetonitrile, Dimethylformamide, and Dimethylsulfoxide at T = (288.15 to 308.15) K |

|         |         |     |        |  |  |
|---------|---------|-----|--------|--|--|
| speedsl | 1457.49 | m/s | 298.15 | Isentropic compressibilities of (amide + water) mixtures: A comparative study  |  |
| speedsl | 1460.20 | m/s | 298.15 | Compressibility Studies of Binary Solutions Involving Water as a Solute in Nonaqueous Solvents at T ) 298.15 K   |  |
| speedsl | 1457.13 | m/s | 298.15 | Thermodynamic Properties of Inorganic Salts in Nonaqueous Solvents. V. Apparent Molar Volumes and Compressibilities of Divalent Transition Metal Bromides in N,N-Dimethylformamide |  |
| srf     | 0.03    | N/m | 327.88 | Surface Tension of Pure Liquids and Binary Liquid Mixtures   |  |
| srf     | 0.04    | N/m | 278.15 | Thermodynamic investigation of N,N-dimethylformamide/toluene binary mixtures in the temperature range from 278.15 to 293.15 K  |  |
| srf     | 0.04    | N/m | 283.15 | Thermodynamic investigation of N,N-dimethylformamide/toluene binary mixtures in the temperature range from 278.15 to 293.15 K  |  |
| srf     | 0.04    | N/m | 293.15 | Thermodynamic investigation of N,N-dimethylformamide/toluene binary mixtures in the temperature range from 278.15 to 293.15 K  |  |
| srf     | 0.04    | N/m | 277.85 | Surface Tension of Pure Liquids and Binary Liquid Mixtures   |  |
| srf     | 0.04    | N/m | 287.81 | Surface Tension of Pure Liquids and Binary Liquid Mixtures   |  |

|     |      |     |        |  |
|-----|------|-----|--------|--|
| srf | 0.04 | N/m | 297.82 | Surface Tension of Pure Liquids and Binary Liquid Mixtures |
| srf | 0.03 | N/m | 307.86 | Surface Tension of Pure Liquids and Binary Liquid Mixtures |
| srf | 0.03 | N/m | 317.86 | Surface Tension of Pure Liquids and Binary Liquid Mixtures |
| srf | 0.03 | N/m | 327.89 | Surface Tension of Pure Liquids and Binary Liquid Mixtures |

Correlations

| Information                 | Value                         |
|-----------------------------|-------------------------------|
| Property code               | pvap                          |
| Equation                    | $\ln(P_{vp}) = A + B/(T + C)$ |
| Coeff. A                    | 1.46587e+01                   |
| Coeff. B                    | -3.67936e+03                  |
| Coeff. C                    | -5.96930e+01                  |
| Temperature range (K), min. | 315.72                        |
| Temperature range (K), max. | 649.60                        |

| Information                 | Value                                      |
|-----------------------------|--|
| Property code               | pvap                                       |
| Equation                    | $\ln(P_{vp}) = A + B/T + C*\ln(T) + D*T^2$ |
| Coeff. A                    | 1.59178e+02                                |
| Coeff. B                    | -1.14422e+04                               |
| Coeff. C                    | -2.15502e+01                               |
| Coeff. D                    | 1.49998e-05                                |
| Temperature range (K), min. | 212.72                                     |
| Temperature range (K), max. | 647.00                                     |

Datasets

## Mass density, kg/m3

| Temperature, K - Liquid | Pressure, kPa - Liquid | Mass density, kg/m3 - Liquid  |
|-------------------------|------------------------|---|
| 298.15                  | 85.90                  | 944.53  |
| Reference               |                        | <a href="https://www.doi.org/10.1016/j.fluid.2013.05.022">https://www.doi.org/10.1016/j.fluid.2013.05.022</a> |

| Temperature, K | Pressure, kPa | Mass density, kg/m3 |
|----------------|---------------|---------------------|
| 313.12         | 1008.00       | 930.41              |
| 313.12         | 2010.00       | 931.09              |
| 313.12         | 3012.00       | 931.73              |
| 313.12         | 4008.00       | 932.38              |
| 313.12         | 5004.00       | 933.04              |
| 313.12         | 6002.00       | 933.68              |
| 313.12         | 7002.00       | 934.32              |
| 313.12         | 8004.00       | 934.97              |
| 313.12         | 9000.00       | 935.62              |
| 313.12         | 10004.00      | 936.27              |
| 313.12         | 11004.00      | 936.89              |
| 313.12         | 12004.00      | 937.51              |
| 313.12         | 13005.00      | 938.14              |
| 313.12         | 14007.00      | 938.78              |
| 313.12         | 15002.00      | 939.37              |
| 313.12         | 16012.00      | 940.0               |
| 313.12         | 17009.00      | 940.59              |
| 313.12         | 18015.00      | 941.19              |
| 313.12         | 19010.00      | 941.79              |
| 313.12         | 20013.00      | 942.42              |
| 313.12         | 21009.00      | 942.99              |
| 313.12         | 22011.00      | 943.58              |
| 313.12         | 23019.00      | 944.17              |
| 313.12         | 24013.00      | 944.75              |
| 313.12         | 25011.00      | 945.35              |
| 323.06         | 1014.00       | 920.89              |
| 323.06         | 2007.00       | 921.59              |
| 323.06         | 3026.00       | 922.27              |
| 323.06         | 4006.00       | 922.95              |
| 323.06         | 5016.00       | 923.62              |
| 323.06         | 6010.00       | 924.3               |
| 323.06         | 7006.00       | 924.99              |
| 323.06         | 8010.00       | 925.64              |

|        |          |        |
|--------|----------|--------|
| 323.06 | 9012.00  | 926.32 |
| 323.06 | 10014.00 | 926.99 |
| 323.06 | 11011.00 | 927.65 |
| 323.06 | 12013.00 | 928.31 |
| 323.06 | 13007.00 | 928.97 |
| 323.06 | 14011.00 | 929.62 |
| 323.06 | 15008.00 | 930.29 |
| 323.06 | 16014.00 | 930.92 |
| 323.06 | 17004.00 | 931.54 |
| 323.06 | 18006.00 | 932.19 |
| 323.06 | 19006.00 | 932.82 |
| 323.06 | 20005.00 | 933.44 |
| 323.06 | 21002.00 | 934.07 |
| 323.06 | 22013.00 | 934.7  |
| 323.06 | 23015.00 | 935.31 |
| 323.06 | 24008.00 | 935.92 |
| 323.06 | 25012.00 | 936.53 |
| 333.00 | 1008.00  | 911.34 |
| 333.00 | 2018.00  | 912.08 |
| 333.00 | 3006.00  | 912.78 |
| 333.00 | 4005.00  | 913.51 |
| 333.00 | 5005.00  | 914.23 |
| 333.00 | 6008.00  | 914.97 |
| 333.00 | 7003.00  | 915.68 |
| 333.00 | 8009.00  | 916.39 |
| 333.00 | 9010.00  | 917.11 |
| 333.00 | 10008.00 | 917.81 |
| 333.00 | 11008.00 | 918.51 |
| 333.00 | 12008.00 | 919.2  |
| 333.00 | 13013.00 | 919.9  |
| 333.00 | 14006.00 | 920.57 |
| 333.00 | 15007.00 | 921.24 |
| 333.00 | 16011.00 | 921.92 |
| 333.00 | 17019.00 | 922.59 |
| 333.00 | 18010.00 | 923.27 |
| 333.00 | 19004.00 | 923.91 |
| 333.00 | 20012.00 | 924.58 |
| 333.00 | 21005.00 | 925.24 |
| 333.00 | 22005.00 | 925.88 |
| 333.00 | 23014.00 | 926.52 |
| 333.00 | 24009.00 | 927.19 |
| 333.00 | 25009.00 | 927.83 |
| 342.86 | 1028.00  | 901.52 |
| 342.86 | 2020.00  | 902.3  |

|        |          |        |
|--------|----------|--------|
| 342.86 | 3015.00  | 903.09 |
| 342.86 | 4002.00  | 903.84 |
| 342.86 | 5035.00  | 904.63 |
| 342.86 | 6026.00  | 905.39 |
| 342.86 | 7013.00  | 906.13 |
| 342.86 | 8017.00  | 906.89 |
| 342.86 | 9024.00  | 907.66 |
| 342.86 | 10012.00 | 908.38 |
| 342.86 | 11048.00 | 909.14 |
| 342.86 | 11999.00 | 909.83 |
| 342.86 | 13018.00 | 910.6  |
| 342.86 | 14029.00 | 911.32 |
| 342.86 | 15007.00 | 912.03 |
| 342.86 | 15994.00 | 912.71 |
| 342.86 | 17041.00 | 913.47 |
| 342.86 | 18012.00 | 914.16 |
| 342.86 | 19006.00 | 914.85 |
| 342.86 | 20042.00 | 915.55 |
| 342.86 | 21025.00 | 916.22 |
| 342.86 | 22008.00 | 916.9  |
| 342.86 | 23079.00 | 917.63 |
| 342.86 | 24008.00 | 918.25 |
| 342.86 | 25017.00 | 918.93 |
| 352.80 | 1040.00  | 891.76 |
| 352.80 | 2009.00  | 892.57 |
| 352.80 | 3011.00  | 893.41 |
| 352.80 | 4020.00  | 894.22 |
| 352.80 | 5050.00  | 895.06 |
| 352.80 | 6060.00  | 895.78 |
| 352.80 | 6996.00  | 896.53 |
| 352.80 | 8008.00  | 897.34 |
| 352.80 | 9034.00  | 898.16 |
| 352.80 | 10019.00 | 898.93 |
| 352.80 | 11043.00 | 899.73 |
| 352.80 | 11998.00 | 900.45 |
| 352.80 | 13028.00 | 901.24 |
| 352.80 | 14015.00 | 901.99 |
| 352.80 | 15028.00 | 902.74 |
| 352.80 | 16000.00 | 903.48 |
| 352.80 | 17011.00 | 904.23 |
| 352.80 | 17999.00 | 904.96 |
| 352.80 | 19044.00 | 905.73 |
| 352.80 | 20050.00 | 906.43 |
| 352.80 | 21038.00 | 907.16 |

|        |          |        |
|--------|----------|--------|
| 352.80 | 21955.00 | 907.8  |
| 352.80 | 23009.00 | 908.57 |
| 352.80 | 23996.00 | 909.27 |
| 352.80 | 25056.00 | 910.04 |
| 362.67 | 1058.00  | 882.04 |
| 362.67 | 2038.00  | 882.89 |
| 362.67 | 3044.00  | 883.77 |
| 362.67 | 4027.00  | 884.62 |
| 362.67 | 5068.00  | 885.54 |
| 362.67 | 6036.00  | 886.37 |
| 362.67 | 7025.00  | 887.22 |
| 362.67 | 8027.00  | 888.03 |
| 362.67 | 9045.00  | 888.9  |
| 362.67 | 10021.00 | 889.71 |
| 362.67 | 11024.00 | 890.55 |
| 362.67 | 11996.00 | 891.3  |
| 362.67 | 13061.00 | 892.19 |
| 362.67 | 14027.00 | 892.95 |
| 362.67 | 15025.00 | 893.73 |
| 362.67 | 15987.00 | 894.49 |
| 362.67 | 17050.00 | 895.32 |
| 362.67 | 18009.00 | 896.08 |
| 362.67 | 19041.00 | 896.85 |
| 362.67 | 20032.00 | 897.62 |
| 362.67 | 21035.00 | 898.36 |
| 362.67 | 22014.00 | 899.1  |
| 362.67 | 23018.00 | 899.87 |
| 362.67 | 24007.00 | 900.58 |
| 362.67 | 25016.00 | 901.35 |

Reference

<https://www.doi.org/10.1021/je050050p>

| Temperature, K | Pressure, kPa | Mass density, kg/m <sup>3</sup> |
|----------------|---------------|---------------------------------|
| 288.15         | 100.00        | 953.4                           |
| 288.15         | 2100.00       | 954.5                           |
| 288.15         | 4000.00       | 955.6                           |
| 288.15         | 6100.00       | 956.8                           |
| 288.15         | 7900.00       | 957.8                           |
| 288.15         | 9900.00       | 958.9                           |
| 288.15         | 12100.00      | 960.1                           |
| 288.15         | 14000.00      | 961.2                           |
| 288.15         | 16000.00      | 962.2                           |
| 288.15         | 18000.00      | 963.3                           |

|        |          |       |
|--------|----------|-------|
| 288.15 | 19800.00 | 964.3 |
| 298.15 | 100.00   | 943.8 |
| 298.15 | 2000.00  | 945.0 |
| 298.15 | 4100.00  | 946.2 |
| 298.15 | 6100.00  | 947.4 |
| 298.15 | 8100.00  | 948.6 |
| 298.15 | 10100.00 | 949.8 |
| 298.15 | 12100.00 | 951.0 |
| 298.15 | 14100.00 | 952.1 |
| 298.15 | 16300.00 | 953.3 |
| 298.15 | 18300.00 | 954.3 |
| 298.15 | 20100.00 | 955.5 |
| 308.15 | 100.00   | 935.0 |
| 308.15 | 2000.00  | 936.2 |
| 308.15 | 4100.00  | 937.6 |
| 308.15 | 6100.00  | 938.9 |
| 308.15 | 8100.00  | 940.0 |
| 308.15 | 10100.00 | 941.2 |
| 308.15 | 12100.00 | 942.5 |
| 308.15 | 14300.00 | 943.7 |
| 308.15 | 16300.00 | 944.9 |
| 308.15 | 18300.00 | 946.0 |
| 308.15 | 20200.00 | 947.0 |
| 318.15 | 100.00   | 925.5 |
| 318.15 | 2100.00  | 926.9 |
| 318.15 | 4000.00  | 928.3 |
| 318.15 | 6000.00  | 929.6 |
| 318.15 | 7900.00  | 930.8 |
| 318.15 | 10200.00 | 932.2 |
| 318.15 | 12000.00 | 933.4 |
| 318.15 | 14000.00 | 934.5 |
| 318.15 | 16000.00 | 935.8 |
| 318.15 | 18000.00 | 937.0 |
| 318.15 | 20200.00 | 938.1 |

Reference

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Sources

Solubility of Etoricoxib in Aqueous Solutions of 1,4-Butanediol, Equilibrium Solubility, Diffusion Coefficient, and Thermodynamic Properties of Etoricoxib in Binary and Ternary Systems of 1,4-Butanediol + Nicotinamide and Nicotinamide in Dimethylsulfoxide + Water) and (Dimethylsulfoxide + Water) Mixed Solvents at 298.15 K:

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|                 |   |
|-----------------|---|
| <b>affp:</b>    | Proton affinity   |
| <b>basg:</b>    | Gas basicity  |
| <b>chl:</b>     | Standard liquid enthalpy of combustion                    |
| <b>cpg:</b>     | Ideal gas heat capacity                                   |
| <b>cpl:</b>     | Liquid phase heat capacity                                |
| <b>dvisc:</b>   | Dynamic viscosity   |
| <b>ea:</b>      | Electron affinity   |
| <b>econd:</b>   | Electrical conductivity                                   |
| <b>gf:</b>      | Standard Gibbs free energy of formation                   |
| <b>hf:</b>      | Enthalpy of formation at standard conditions              |
| <b>hfl:</b>     | Liquid phase enthalpy of formation at standard conditions |
| <b>hfus:</b>    | Enthalpy of fusion at standard conditions                 |
| <b>hfust:</b>   | Enthalpy of fusion at a given temperature                 |
| <b>hvap:</b>    | Enthalpy of vaporization at standard conditions           |
| <b>hvapt:</b>   | Enthalpy of vaporization at a given temperature           |
| <b>ie:</b>      | Ionization energy   |
| <b>log10ws:</b> | Log10 of Water solubility in mol/l                        |
| <b>logp:</b>    | Octanol/Water partition coefficient                       |
| <b>mcvol:</b>   | McGowan's characteristic volume                           |
| <b>nfpaf:</b>   | NFPA Fire Rating  |
| <b>nfpah:</b>   | NFPA Health Rating  |
| <b>pc:</b>      | Critical Pressure   |
| <b>pvap:</b>    | Vapor pressure  |
| <b>rfi:</b>     | Refractive Index  |
| <b>rhoc:</b>    | Critical density  |
| <b>rho:</b>     | Liquid Density  |
| <b>rinpol:</b>  | Non-polar retention indices                               |
| <b>ripol:</b>   | Polar retention indices                                   |
| <b>sfust:</b>   | Entropy of fusion at a given temperature                  |
| <b>speedsl:</b> | Speed of sound in fluid                                   |
| <b>srf:</b>     | Surface Tension   |
| <b>tb:</b>      | Normal Boiling Point Temperature                          |
| <b>tc:</b>      | Critical Temperature                                      |
| <b>tf:</b>      | Normal melting (fusion) point                             |
| <b>vc:</b>      | Critical Volume   |

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