

# Tetrahydro-4H-pyran-4-ol

|                             |  |
|-----------------------------|--|
| <b>Other names:</b>         | 2H-Pyran-4-ol, tetrahydro-tetrahydro-2H-pyran-4-ol |
| <b>Inchi:</b>               | InChI=1S/C5H10O2/c6-5-1-3-7-4-2-5/h5-6H,1-4H2      |
| <b>InchiKey:</b>            | LMYJGUNNJIDROI-UHFFFAOYSA-N                        |
| <b>Formula:</b>             | C5H10O2  |
| <b>SMILES:</b>              | OC1CCOCC1  |
| <b>Mol. weight [g/mol]:</b> | 102.13   |
| <b>CAS:</b>                 | 2081-44-9  |

## Physical Properties

| Property code | Value   | Unit                 | Source         |
|---------------|---------|----------------------|----------------|
| gf            | -207.27 | kJ/mol               | Joback Method  |
| hf            | -376.44 | kJ/mol               | Joback Method  |
| hfus          | 12.61   | kJ/mol               | Joback Method  |
| hvap          | 48.34   | kJ/mol               | Joback Method  |
| log10ws       | -0.27   |                      | Crippen Method |
| logp          | 0.158   |                      | Crippen Method |
| mcvol         | 82.190  | ml/mol               | McGowan Method |
| pc            | 5029.93 | kPa                  | Joback Method  |
| tb            | 452.48  | K                    | Joback Method  |
| tc            | 649.36  | K                    | Joback Method  |
| tf            | 240.88  | K                    | Joback Method  |
| vc            | 0.288   | m <sup>3</sup> /kmol | Joback Method  |

## Temperature Dependent Properties

| Property code | Value  | Unit    | Temperature [K] | Source        |
|---------------|--------|---------|-----------------|---------------|
| cpg           | 175.96 | J/molxK | 452.48          | Joback Method |
| cpg           | 227.36 | J/molxK | 616.54          | Joback Method |
| cpg           | 218.15 | J/molxK | 583.73          | Joback Method |
| cpg           | 208.42 | J/molxK | 550.92          | Joback Method |
| cpg           | 198.15 | J/molxK | 518.11          | Joback Method |
| cpg           | 187.33 | J/molxK | 485.29          | Joback Method |
| cpg           | 236.06 | J/molxK | 649.36          | Joback Method |

|       |           |      |        |               |
|-------|-----------|------|--------|---------------|
| dvisc | 0.0002682 | Paxs | 452.48 | Joback Method |
| dvisc | 0.0004569 | Paxs | 417.21 | Joback Method |
| dvisc | 0.0008590 | Paxs | 381.95 | Joback Method |
| dvisc | 0.0018359 | Paxs | 346.68 | Joback Method |
| dvisc | 0.0046608 | Paxs | 311.41 | Joback Method |
| dvisc | 0.0150108 | Paxs | 276.15 | Joback Method |
| dvisc | 0.0680898 | Paxs | 240.88 | Joback Method |

## Pressure Dependent Properties

| Property code | Value  | Unit | Pressure [kPa] | Source       |
|---------------|--------|------|----------------|--------------|
| tbrp          | 360.20 | K    | 2.00           | NIST Webbook |

## Sources

|                        |   |
|------------------------|---|
| <b>Joback Method:</b>  | <a href="https://en.wikipedia.org/wiki/Joback_method">https://en.wikipedia.org/wiki/Joback_method</a>                                       |
| <b>McGowan Method:</b> | <a href="http://link.springer.com/article/10.1007/BF02311772">http://link.springer.com/article/10.1007/BF02311772</a>                       |
| <b>NIST Webbook:</b>   | <a href="http://webbook.nist.gov/cgi/cbook.cgi?ID=C2081449&amp;Units=SI">http://webbook.nist.gov/cgi/cbook.cgi?ID=C2081449&amp;Units=SI</a> |
| <b>Crippen Method:</b> | <a href="http://pubs.acs.org/doi/abs/10.1021/ci990307l">http://pubs.acs.org/doi/abs/10.1021/ci990307l</a>                                   |
| <b>Crippen Method:</b> | <a href="https://www.chemeo.com/doc/models/crippen_log10ws">https://www.chemeo.com/doc/models/crippen_log10ws</a>                           |

## Legend

|                 |   |
|-----------------|---|
| <b>cpg:</b>     | Ideal gas heat capacity                         |
| <b>dvisc:</b>   | Dynamic viscosity                               |
| <b>gf:</b>      | Standard Gibbs free energy of formation         |
| <b>hf:</b>      | Enthalpy of formation at standard conditions    |
| <b>hfus:</b>    | Enthalpy of fusion at standard conditions       |
| <b>hvap:</b>    | Enthalpy of vaporization at standard conditions |
| <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>pc:</b>      | Critical Pressure                               |
| <b>tb:</b>      | Normal Boiling Point Temperature                |
| <b>tbrp:</b>    | Boiling point at reduced pressure               |

**tc:** Critical Temperature  
**tf:** Normal melting (fusion) point  
**vc:** Critical Volume

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