

# 3(2H)-Thiophenone, dihydro-2-methyl-

|                             |  |
|-----------------------------|--|
| <b>Other names:</b>         | Dihydro-2-methyl-3(2H)-thiophenone<br>2-Methyltetrahydrothiophen-3-one<br>2-Methyldihydro-3(2H)-thiophenone<br>2-Methyl-3-thiolanone<br>2-methyl-4,5-dihydro-3(2H)-thiophenone<br>2-Methyltetrahydrothiophene-3-one<br>2-Methylthiolan-3-one<br>2-Methyltetrahydrothiophen-3-one<br>4,5-Dihydro-2-methyl-3(2H)-thiophenone<br>4,5-Dihydro-2-methylthiophen-3(2H)-one<br>Dihydro-3(2H)-thiophenone, 2-methyl<br>Thiolan-3-one, 2-methyl<br>Thiophen-3(2H)-one, dihydro-2-methyl<br>dihydro-2-methylthiophen-3(2H)-one |
| <b>Inchi:</b>               | InChI=1S/C5H8OS/c1-4-5(6)2-3-7-4/h4H,2-3H2,1H3   |
| <b>InchiKey:</b>            | YMZZPMVKABUEBL-UHFFFAOYSA-N  |
| <b>Formula:</b>             | C5H8OS   |
| <b>SMILES:</b>              | CC1SCCC1=O   |
| <b>Mol. weight [g/mol]:</b> | 116.18   |
| <b>CAS:</b>                 | 13679-85-1   |

## Physical Properties

| Property code | Value   | Unit   | Source         |
|---------------|---------|--------|----------------|
| gf            | -54.96  | kJ/mol | Joback Method  |
| hf            | -178.49 | kJ/mol | Joback Method  |
| hfus          | 5.81    | kJ/mol | Joback Method  |
| hvap          | 37.04   | kJ/mol | Joback Method  |
| log10ws       | -1.09   |        | Crippen Method |
| logp          | 1.081   |        | Crippen Method |
| mcvol         | 88.370  | ml/mol | McGowan Method |
| pc            | 4510.35 | kPa    | Joback Method  |
| rinpol        | 1017.00 |        | NIST Webbook   |
| rinpol        | 990.00  |        | NIST Webbook   |
| rinpol        | 1012.00 |        | NIST Webbook   |
| rinpol        | 990.00  |        | NIST Webbook   |
| rinpol        | 947.00  |        | NIST Webbook   |
| rinpol        | 952.00  |        | NIST Webbook   |

|        |         |              |
|--------|---------|--------------|
| rinpol | 951.00  | NIST Webbook |
| rinpol | 956.00  | NIST Webbook |
| rinpol | 1000.00 | NIST Webbook |
| rinpol | 1011.00 | NIST Webbook |
| rinpol | 1009.00 | NIST Webbook |
| rinpol | 1017.00 | NIST Webbook |
| rinpol | 996.00  | NIST Webbook |
| rinpol | 1017.00 | NIST Webbook |
| rinpol | 989.00  | NIST Webbook |
| rinpol | 958.00  | NIST Webbook |
| rinpol | 952.00  | NIST Webbook |
| rinpol | 947.00  | NIST Webbook |
| rinpol | 1001.00 | NIST Webbook |
| rinpol | 994.00  | NIST Webbook |
| rinpol | 996.00  | NIST Webbook |
| rinpol | 949.00  | NIST Webbook |
| rinpol | 947.00  | NIST Webbook |
| rinpol | 1000.00 | NIST Webbook |
| rinpol | 1017.00 | NIST Webbook |
| rinpol | 958.00  | NIST Webbook |
| rinpol | 1001.00 | NIST Webbook |
| rinpol | 999.00  | NIST Webbook |
| rinpol | 998.00  | NIST Webbook |
| rinpol | 990.00  | NIST Webbook |
| rinpol | 979.00  | NIST Webbook |
| rinpol | 994.00  | NIST Webbook |
| rinpol | 990.00  | NIST Webbook |
| ripol  | 1518.00 | NIST Webbook |
| ripol  | 1506.00 | NIST Webbook |
| ripol  | 1561.00 | NIST Webbook |
| ripol  | 1551.00 | NIST Webbook |
| ripol  | 1557.00 | NIST Webbook |
| ripol  | 1551.00 | NIST Webbook |
| ripol  | 1551.00 | NIST Webbook |
| ripol  | 1518.00 | NIST Webbook |
| ripol  | 1520.00 | NIST Webbook |
| ripol  | 1506.00 | NIST Webbook |
| ripol  | 1511.00 | NIST Webbook |
| ripol  | 1520.00 | NIST Webbook |
| ripol  | 1512.00 | NIST Webbook |
| ripol  | 1512.00 | NIST Webbook |
| ripol  | 1535.00 | NIST Webbook |
| ripol  | 1533.00 | NIST Webbook |
| ripol  | 1548.00 | NIST Webbook |

|       |         |                      |               |
|-------|---------|----------------------|---------------|
| ripol | 1565.00 |                      | NIST Webbook  |
| ripol | 1538.00 |                      | NIST Webbook  |
| ripol | 1538.00 |                      | NIST Webbook  |
| ripol | 1565.00 |                      | NIST Webbook  |
| ripol | 1542.00 |                      | NIST Webbook  |
| ripol | 1509.00 |                      | NIST Webbook  |
| ripol | 1518.00 |                      | NIST Webbook  |
| ripol | 1528.00 |                      | NIST Webbook  |
| ripol | 1524.00 |                      | NIST Webbook  |
| ripol | 1535.00 |                      | NIST Webbook  |
| ripol | 1538.00 |                      | NIST Webbook  |
| ripol | 1523.00 |                      | NIST Webbook  |
| ripol | 1506.00 |                      | NIST Webbook  |
| ripol | 1512.00 |                      | NIST Webbook  |
| ripol | 1510.00 |                      | NIST Webbook  |
| ripol | 1506.00 |                      | NIST Webbook  |
| ripol | 1525.00 |                      | NIST Webbook  |
| ripol | 1510.00 |                      | NIST Webbook  |
| tb    | 444.73  | K                    | Joback Method |
| tc    | 682.05  | K                    | Joback Method |
| tf    | 308.68  | K                    | Joback Method |
| vc    | 0.309   | m <sup>3</sup> /kmol | Joback Method |

## Temperature Dependent Properties

| Property code | Value  | Unit    | Temperature [K] | Source        |
|---------------|--------|---------|-----------------|---------------|
| cpg           | 168.02 | J/mol×K | 444.73          | Joback Method |
| cpg           | 180.14 | J/mol×K | 484.28          | Joback Method |
| cpg           | 191.71 | J/mol×K | 523.84          | Joback Method |
| cpg           | 202.73 | J/mol×K | 563.39          | Joback Method |
| cpg           | 213.20 | J/mol×K | 602.94          | Joback Method |
| cpg           | 223.10 | J/mol×K | 642.49          | Joback Method |
| cpg           | 232.43 | J/mol×K | 682.05          | Joback Method |

## Sources

McGowan Method:

<http://link.springer.com/article/10.1007/BF02311772>

NIST Webbook:

<http://webbook.nist.gov/cgi/cbook.cgi?ID=C13679851&Units=SI>

**Crippen Method:** <http://pubs.acs.org/doi/abs/10.1021/ci9903071>  
**Crippen Method:** [https://www.chemeo.com/doc/models/crippen\\_log10ws](https://www.chemeo.com/doc/models/crippen_log10ws)  
**Joback Method:** [https://en.wikipedia.org/wiki/Joback\\_method](https://en.wikipedia.org/wiki/Joback_method)

## Legend

**cpg:** Ideal gas heat capacity  
**gf:** Standard Gibbs free energy of formation  
**hf:** Enthalpy of formation at standard conditions  
**hfus:** Enthalpy of fusion at standard conditions  
**hvap:** Enthalpy of vaporization at standard conditions  
**log10ws:** Log10 of Water solubility in mol/l  
**logp:** Octanol/Water partition coefficient  
**mcvol:** McGowan's characteristic volume  
**pc:** Critical Pressure  
**rinpol:** Non-polar retention indices  
**ripol:** Polar retention indices  
**tb:** Normal Boiling Point Temperature  
**tc:** Critical Temperature  
**tf:** Normal melting (fusion) point  
**vc:** Critical Volume

Latest version available from:

<https://www.chemeo.com/cid/25-592-7/3-2H-Thiophenone-dihydro-2-methyl.pdf>

Generated by Cheméo on 2024-09-27 17:26:11.709314653 +0000 UTC m=+2038834.346283901.

Cheméo (<https://www.chemeo.com>) is the biggest free database of chemical and physical data for the process industry.