

2H-Thiopyran, tetrahydro-2-methyl-

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| Other names: | Tetrahydro-2-methyl-2H-thiopyran 2-Methylthiane 2-methyl-thiacyclohexane 2-Methyltetrahydro-2H-thiopyrane |
| Inchi: | InChI=1S/C6H12S/c1-6-4-2-3-5-7-6/h6H,2-5H2,1H3 |
| InchiKey: | VLBGYQISAJHVAD-UHFFFAOYSA-N |
| Formula: | C6H12S |
| SMILES: | CC1CCCCS1 |
| Mol. weight [g/mol]: | 116.22 |
| CAS: | 5161-16-0 |

Physical Properties

| Property code | Value | Unit | Source |
|---------------|---------|---------|----------------|
| gf | 63.95 | kJ/mol | Joback Method |
| hf | -67.59 | kJ/mol | Joback Method |
| hfus | 6.79 | kJ/mol | Joback Method |
| hvap | 35.19 | kJ/mol | Joback Method |
| log10ws | -2.22 | | Crippen Method |
| logp | 2.292 | | Crippen Method |
| mcvol | 100.890 | ml/mol | McGowan Method |
| pc | 3930.78 | kPa | Joback Method |
| rinpol | 924.00 | | NIST Webbook |
| rinpol | 910.00 | | NIST Webbook |
| rinpol | 924.00 | | NIST Webbook |
| rinpol | 902.00 | | NIST Webbook |
| rinpol | 902.00 | | NIST Webbook |
| rinpol | 902.00 | | NIST Webbook |
| rinpol | 924.00 | | NIST Webbook |
| rinpol | 901.00 | | NIST Webbook |
| tb | 404.06 | K | Joback Method |
| tc | 626.29 | K | Joback Method |
| tf | 248.21 | K | Joback Method |
| vc | 0.350 | m3/kmol | Joback Method |

Temperature Dependent Properties

| Property code | Value | Unit | Temperature [K] | Source |
|---------------|--------|---------|-----------------|---------------|
| cpg | 236.46 | J/mol×K | 552.22 | Joback Method |
| cpg | 248.61 | J/mol×K | 589.25 | Joback Method |
| cpg | 180.17 | J/mol×K | 404.06 | Joback Method |
| cpg | 195.45 | J/mol×K | 441.10 | Joback Method |
| cpg | 209.91 | J/mol×K | 478.14 | Joback Method |
| cpg | 223.57 | J/mol×K | 515.18 | Joback Method |
| cpg | 260.03 | J/mol×K | 626.29 | Joback Method |
| hvapt | 42.10 | kJ/mol | 386.00 | NIST Webbook |
| hvapt | 40.20 | kJ/mol | 397.00 | NIST Webbook |

Sources

| | |
|------------------------|---|
| Crippen Method: | https://www.chemeo.com/doc/models/crippen_log10ws |
| Joback Method: | https://en.wikipedia.org/wiki/Joback_method |
| McGowan Method: | http://link.springer.com/article/10.1007/BF02311772 |
| NIST Webbook: | http://webbook.nist.gov/cgi/cbook.cgi?ID=C5161160&Units=SI |
| Crippen Method: | http://pubs.acs.org/doi/abs/10.1021/ci9903071 |

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 |
| hvapt: | Enthalpy of vaporization at a given temperature |
| 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 |
| tb: | Normal Boiling Point Temperature |
| tc: | Critical Temperature |

tf: Normal melting (fusion) point

vc: Critical Volume

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