

CH₃C(O)OCH(CH₂CH₃)C(CH₃)₃

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|-----------------------------|---|
| Other names: | Acetic acid, 2,2-dimethylpent-3-yl ester |
| Inchi: | InChI=1S/C9H18O2/c1-6-8(9(3,4)5)11-7(2)10/h8H,6H2,1-5H3 |
| InchiKey: | OZVHACQHIBUHOC-UHFFFAOYSA-N |
| Formula: | C ₉ H ₁₈ O ₂ |
| SMILES: | CCC(OC(C)=O)C(C)(C)C |
| Mol. weight [g/mol]: | 158.24 |
| CAS: | 39511-81-4 |

Physical Properties

| Property code | Value | Unit | Source |
|---------------|---------|----------------------|----------------|
| gf | -208.62 | kJ/mol | Joback Method |
| hf | -487.92 | kJ/mol | Joback Method |
| hfus | 10.92 | kJ/mol | Joback Method |
| hvap | 43.10 | kJ/mol | Joback Method |
| log10ws | -2.32 | | Crippen Method |
| logp | 2.374 | | Crippen Method |
| mcvol | 145.110 | ml/mol | McGowan Method |
| pc | 2477.65 | kPa | Joback Method |
| rinpol | 970.00 | | NIST Webbook |
| tb | 477.94 | K | Joback Method |
| tc | 666.16 | K | Joback Method |
| tf | 250.77 | K | Joback Method |
| vc | 0.546 | m ³ /kmol | Joback Method |

Temperature Dependent Properties

| Property code | Value | Unit | Temperature [K] | Source |
|---------------|--------|---------|-----------------|---------------|
| cpg | 325.82 | J/mol×K | 477.94 | Joback Method |
| cpg | 392.85 | J/mol×K | 634.79 | Joback Method |
| cpg | 380.76 | J/mol×K | 603.42 | Joback Method |
| cpg | 368.03 | J/mol×K | 572.05 | Joback Method |
| cpg | 354.64 | J/mol×K | 540.68 | Joback Method |
| cpg | 340.58 | J/mol×K | 509.31 | Joback Method |
| cpg | 404.33 | J/mol×K | 666.16 | Joback Method |

| | | | | |
|-------|-----------|------|--------|---------------|
| dvisc | 0.0002253 | Paxs | 477.94 | Joback Method |
| dvisc | 0.0003117 | Paxs | 440.08 | Joback Method |
| dvisc | 0.0004585 | Paxs | 402.22 | Joback Method |
| dvisc | 0.0007309 | Paxs | 364.36 | Joback Method |
| dvisc | 0.0012980 | Paxs | 326.49 | Joback Method |
| dvisc | 0.0026801 | Paxs | 288.63 | Joback Method |
| dvisc | 0.0068881 | Paxs | 250.77 | Joback Method |

Sources

| | |
|------------------------|---|
| Crippen Method: | http://pubs.acs.org/doi/abs/10.1021/ci990307l |
| 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=C39511814&Units=SI |

Legend

| | |
|-----------------|---|
| cpg: | Ideal gas heat capacity |
| dvisc: | Dynamic viscosity |
| 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 |
| tb: | Normal Boiling Point Temperature |
| tc: | Critical Temperature |
| tf: | Normal melting (fusion) point |
| vc: | Critical Volume |

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<https://www.chemeo.com/cid/47-058-6/CH3C-O-OCH-CH2CH3-C-CH3-3.pdf>

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