

2-Cyclohexen-1-one, 3-methyl-6-(1-methylethyl)-

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| Other names: | p-Menth-1-en-3-one Piperitone 3-Carvomenthene 1-Methyl-4-isopropyl-1-cyclohexen-3-one 3-Methyl-6-(1-methylethyl)-2-cyclohexen-1-one 3-Methyl-6-isopropyl-2-cyclohexen-1-one 6-isopropyl-3-methylcyclohex-2-enone |
| Inchi: | InChI=1S/C10H16O/c1-7(2)9-5-4-8(3)6-10(9)11/h6-7,9H,4-5H2,1-3H3 |
| InchiKey: | YSTPAHQEHQSRJD-UHFFFAOYSA-N |
| Formula: | C10H16O |
| SMILES: | CC1=CC(=O)C(C(C)C)CC1 |
| Mol. weight [g/mol]: | 152.23 |
| CAS: | 89-81-6 |

Physical Properties

| Property code | Value | Unit | Source |
|---------------|---------|--------|----------------|
| gf | -46.93 | kJ/mol | Joback Method |
| hf | -292.08 | kJ/mol | Joback Method |
| hfus | 10.31 | kJ/mol | Joback Method |
| hvap | 43.10 | kJ/mol | Joback Method |
| log10ws | -2.55 | | Crippen Method |
| logp | 2.568 | | Crippen Method |
| mcvol | 138.170 | ml/mol | McGowan Method |
| pc | 2744.03 | kPa | Joback Method |
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| tc | 741.04 | K | Joback Method |
| tf | 276.34 | K | Joback Method |
| vc | 0.515 | m ³ /kmol | Joback Method |

Temperature Dependent Properties

| Property code | Value | Unit | Temperature [K] | Source |
|---------------|--------|---------|-----------------|---------------|
| cpg | 321.40 | J/mol×K | 519.27 | Joback Method |
| cpg | 339.21 | J/mol×K | 556.23 | Joback Method |
| cpg | 356.17 | J/mol×K | 593.19 | Joback Method |
| cpg | 372.28 | J/mol×K | 630.16 | Joback Method |
| cpg | 387.53 | J/mol×K | 667.12 | Joback Method |
| cpg | 401.93 | J/mol×K | 704.08 | Joback Method |
| cpg | 415.47 | J/mol×K | 741.04 | Joback Method |

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Crippen Method:

<http://pubs.acs.org/doi/abs/10.1021/ci9903071>

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>

Legend

| | |
|---------------------------------------|---|
| cp_g: | Ideal gas heat capacity |
| g_f: | Standard Gibbs free energy of formation |
| h_f: | Enthalpy of formation at standard conditions |
| h_{fus}: | Enthalpy of fusion at standard conditions |
| h_{vap}: | Enthalpy of vaporization at standard conditions |
| log₁₀w_s: | Log ₁₀ of Water solubility in mol/l |
| log_p: | Octanol/Water partition coefficient |
| mc_{vol}: | McGowan's characteristic volume |
| pc: | Critical Pressure |
| rin_{pol}: | Non-polar retention indices |
| rip_{ol}: | Polar retention indices |
| tb: | Normal Boiling Point Temperature |
| tc: | Critical Temperature |
| tf: | Normal melting (fusion) point |
| vc: | Critical Volume |

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