

Cyclohexane, 1-methyl-4-(1-methylethylidene)-

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| Other names: | p-Menth-4(8)-ene 4-Methyl-1-isopropylidenecyclohexane 1-Isopropylidene-4-methylcyclohexane 4(8)-p-Menthene «DELTA»4(8)-Menthene p-Ment-4(8)-ene |
| Inchi: | InChI=1S/C10H18/c1-8(2)10-6-4-9(3)5-7-10/h9H,4-7H2,1-3H3 |
| InchiKey: | OLZFGHOLNXRJEC-UHFFFAOYSA-N |
| Formula: | C10H18 |
| SMILES: | CC(C)=C1CCC(C)CC1 |
| Mol. weight [g/mol]: | 138.25 |
| CAS: | 1124-27-2 |

Physical Properties

| Property code | Value | Unit | Source |
|---------------|---------------|----------------------|----------------|
| gf | 94.68 | kJ/mol | Joback Method |
| hf | -129.17 | kJ/mol | Joback Method |
| hfus | 12.50 | kJ/mol | Joback Method |
| hvap | 39.15 | kJ/mol | Joback Method |
| log10ws | -3.52 | | Crippen Method |
| logp | 3.533 | | Crippen Method |
| mcvol | 136.600 | ml/mol | McGowan Method |
| pc | 2659.77 | kPa | Joback Method |
| rinpol | 998.00 | | NIST Webbook |
| rinpol | 1025.90 | | NIST Webbook |
| rinpol | 1018.00 | | NIST Webbook |
| rinpol | 998.00 | | NIST Webbook |
| rinpol | 1025.90 | | NIST Webbook |
| ripol | 1220.00 | | NIST Webbook |
| tb | 446.00 ± 6.00 | K | NIST Webbook |
| tb | 446.00 ± 6.00 | K | NIST Webbook |
| tc | 662.43 | K | Joback Method |
| tf | 206.24 | K | Joback Method |
| vc | 0.512 | m ³ /kmol | Joback Method |

Temperature Dependent Properties

| Property code | Value | Unit | Temperature [K] | Source |
|---------------|--------|---------|-----------------|---------------|
| cpg | 286.34 | J/molxK | 454.27 | Joback Method |
| cpg | 305.04 | J/molxK | 488.96 | Joback Method |
| cpg | 322.79 | J/molxK | 523.66 | Joback Method |
| cpg | 339.62 | J/molxK | 558.35 | Joback Method |
| cpg | 355.55 | J/molxK | 593.04 | Joback Method |
| cpg | 370.62 | J/molxK | 627.74 | Joback Method |
| cpg | 384.85 | J/molxK | 662.43 | Joback Method |

Sources

| | |
|------------------------|---|
| NIST Webbook: | http://webbook.nist.gov/cgi/cbook.cgi?ID=C1124272&Units=SI |
| 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 |

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 |
| ripol: | Non-polar retention indices |
| ripol: | Polar retention indices |
| tb: | Normal Boiling Point Temperature |
| tc: | Critical Temperature |
| tf: | Normal melting (fusion) point |
| vc: | Critical Volume |

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