

Cyclohexane, 1-methyl-4-(1-methylethenyl)-, trans-

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| Other names: | p-Menth-8-ene, trans- trans-p-Menth-8-ene trans-1-Isopropenyl-4-methylcyclohexane 8-p-Menthene (trans) 1-Isopropenyl-4-methylcyclohexane, (E)- |
| Inchi: | InChI=1S/C10H18/c1-8(2)10-6-4-9(3)5-7-10/h9-10H,1,4-7H2,2-3H3/t9-,10- |
| InchiKey: | WPMKLOWQWIDOJN-MGCOHNPYSA-N |
| Formula: | C10H18 |
| SMILES: | C=C(C)C1CCC(C)CC1 |
| Mol. weight [g/mol]: | 138.25 |
| CAS: | 1124-25-0 |

Physical Properties

| Property code | Value | Unit | Source |
|---------------|---------------|----------------------|----------------|
| gf | 129.35 | kJ/mol | Joback Method |
| hf | -100.11 | kJ/mol | Joback Method |
| hfus | 11.97 | kJ/mol | Joback Method |
| hvap | 37.38 | kJ/mol | Joback Method |
| log10ws | -3.27 | | Crippen Method |
| logp | 3.389 | | Crippen Method |
| mcvol | 136.600 | ml/mol | McGowan Method |
| pc | 2605.74 | kPa | Joback Method |
| rinpol | 980.30 | | NIST Webbook |
| rinpol | 974.10 | | NIST Webbook |
| tb | 434.20 ± 2.00 | K | NIST Webbook |
| tb | 516.15 ± 3.00 | K | NIST Webbook |
| tc | 644.45 | K | Joback Method |
| tf | 189.88 | K | Joback Method |
| vc | 0.509 | m ³ /kmol | Joback Method |

Temperature Dependent Properties

| Property code | Value | Unit | Temperature [K] | Source |
|---------------|--------|---------|-----------------|---------------|
| cpg | 282.84 | J/mol×K | 439.64 | Joback Method |

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|-----|--------|---------|--------|---------------|
| cpg | 302.49 | J/mol×K | 473.78 | Joback Method |
| cpg | 321.16 | J/mol×K | 507.91 | Joback Method |
| cpg | 338.90 | J/mol×K | 542.05 | Joback Method |
| cpg | 355.71 | J/mol×K | 576.18 | Joback Method |
| cpg | 371.63 | J/mol×K | 610.32 | Joback Method |
| cpg | 386.68 | J/mol×K | 644.45 | Joback Method |

Sources

| | |
|------------------------|---|
| NIST Webbook: | http://webbook.nist.gov/cgi/cbook.cgi?ID=C1124250&Units=SI |
| 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

| | |
|-----------------|---|
| 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 |
| tb: | Normal Boiling Point Temperature |
| tc: | Critical Temperature |
| tf: | Normal melting (fusion) point |
| vc: | Critical Volume |

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