

Butane, 1-(2,2-dichloro-1-methylcyclopropyl)-3-methyl-

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| Other names: | 1,1-Dichloro-2-isopentyl-2-methylcyclopropane Cyclopropane, 1,1-dichloro-2-methyl-2-(3-methylbutyl) |
| Inchi: | InChI=1S/C9H16Cl2/c1-7(2)4-5-8(3)6-9(8,10)11/h7H,4-6H2,1-3H3 |
| InchiKey: | SZDCIZLZSRDCSY-UHFFFAOYSA-N |
| Formula: | C9H16Cl2 |
| SMILES: | CC(C)CCC1(C)CC1(Cl)Cl |
| Mol. weight [g/mol]: | 195.13 |
| CAS: | 24551-81-3 |

Physical Properties

| Property code | Value | Unit | Source |
|---------------|---------|----------------------|----------------|
| gf | 40.66 | kJ/mol | Joback Method |
| hf | -182.91 | kJ/mol | Joback Method |
| hfus | 10.55 | kJ/mol | Joback Method |
| hvap | 41.31 | kJ/mol | Joback Method |
| log10ws | -3.91 | | Crippen Method |
| logp | 4.006 | | Crippen Method |
| mcvol | 151.290 | ml/mol | McGowan Method |
| pc | 2600.43 | kPa | Joback Method |
| rinpola | 1134.00 | | NIST Webbook |
| ripola | 1364.00 | | NIST Webbook |
| ripolb | 1364.00 | | NIST Webbook |
| tb | 482.29 | K | Joback Method |
| tc | 692.27 | K | Joback Method |
| tf | 297.53 | K | Joback Method |
| vc | 0.584 | m ³ /kmol | Joback Method |

Temperature Dependent Properties

| Property code | Value | Unit | Temperature [K] | Source |
|---------------|--------|---------|-----------------|---------------|
| cpg | 322.58 | J/molxK | 482.29 | Joback Method |
| cpg | 337.72 | J/molxK | 517.29 | Joback Method |
| cpg | 351.56 | J/molxK | 552.28 | Joback Method |
| cpg | 364.32 | J/molxK | 587.28 | Joback Method |

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|-----|--------|---------|--------|---------------|
| cpg | 376.17 | J/mol×K | 622.28 | Joback Method |
| cpg | 387.34 | J/mol×K | 657.28 | Joback Method |
| cpg | 398.01 | J/mol×K | 692.27 | Joback Method |

Sources

| | |
|------------------------|---|
| 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 |
| NIST Webbook: | http://webbook.nist.gov/cgi/cbook.cgi?ID=C24551813&Units=SI |

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

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