

# Cyclohexane, 2-butyl-1,1,3-trimethyl-

|                             |   |
|-----------------------------|---|
| <b>Other names:</b>         | 1,1,3-Trimethyl-2-butyl-cyclohexane<br>2-Butyl-1,1,3-trimethylcyclohexane |
| <b>Inchi:</b>               | InChI=1S/C13H26/c1-5-6-9-12-11(2)8-7-10-13(12,3)4/h11-12H,5-10H2,1-4H3    |
| <b>InchiKey:</b>            | PUUCZCJOATUHNE-UHFFFAOYSA-N   |
| <b>Formula:</b>             | C13H26  |
| <b>SMILES:</b>              | CCCCC1C(C)CCCC1(C)C   |
| <b>Mol. weight [g/mol]:</b> | 182.35  |
| <b>CAS:</b>                 | 54676-39-0  |

## Physical Properties

| Property code | Value   | Unit                 | Source         |
|---------------|---------|----------------------|----------------|
| gf            | 62.12   | kJ/mol               | Joback Method  |
| hf            | -282.77 | kJ/mol               | Joback Method  |
| hfus          | 17.11   | kJ/mol               | Joback Method  |
| hvap          | 43.19   | kJ/mol               | Joback Method  |
| log10ws       | -4.43   |                      | Crippen Method |
| logp          | 4.639   |                      | Crippen Method |
| mvol          | 183.170 | ml/mol               | McGowan Method |
| pc            | 1920.30 | kPa                  | Joback Method  |
| rinpol        | 1228.00 |                      | NIST Webbook   |
| tb            | 507.29  | K                    | Joback Method  |
| tc            | 704.36  | K                    | Joback Method  |
| tf            | 259.07  | K                    | Joback Method  |
| vc            | 0.693   | m <sup>3</sup> /kmol | Joback Method  |

## Temperature Dependent Properties

| Property code | Value  | Unit    | Temperature [K] | Source        |
|---------------|--------|---------|-----------------|---------------|
| cpg           | 445.22 | J/mol×K | 507.29          | Joback Method |
| cpg           | 467.63 | J/mol×K | 540.14          | Joback Method |
| cpg           | 488.86 | J/mol×K | 572.98          | Joback Method |
| cpg           | 509.00 | J/mol×K | 605.83          | Joback Method |
| cpg           | 528.12 | J/mol×K | 638.67          | Joback Method |
| cpg           | 546.32 | J/mol×K | 671.52          | Joback Method |

## Correlations

| Information                 | Value                         |
|-----------------------------|-------------------------------|
| Property code               | pvap                          |
| Equation                    | $\ln(P_{vp}) = A + B/(T + C)$ |
| Coeff. A                    | 1.48084e+01                   |
| Coeff. B                    | -4.28856e+03                  |
| Coeff. C                    | -8.30740e+01                  |
| Temperature range (K), min. | 378.41                        |
| Temperature range (K), max. | 534.65                        |

## Sources

|   |   |
|---|---|
| <b>Crippen Method:</b>                      | <a href="http://pubs.acs.org/doi/abs/10.1021/ci990307l">http://pubs.acs.org/doi/abs/10.1021/ci990307l</a>   |
| <b>Crippen Method:</b>                      | <a href="https://www.chemeo.com/doc/models/crippen_log10ws">https://www.chemeo.com/doc/models/crippen_log10ws</a>   |
| <b>Joback Method:</b>                       | <a href="https://en.wikipedia.org/wiki/Joback_method">https://en.wikipedia.org/wiki/Joback_method</a>   |
| <b>McGowan Method:</b>                      | <a href="http://link.springer.com/article/10.1007/BF02311772">http://link.springer.com/article/10.1007/BF02311772</a>   |
| <b>NIST Webbook:</b>                        | <a href="http://webbook.nist.gov/cgi/cbook.cgi?ID=C54676390&amp;Units=SI">http://webbook.nist.gov/cgi/cbook.cgi?ID=C54676390&amp;Units=SI</a>   |
| <b>The Yaws Handbook of Vapor Pressure:</b> | <a href="https://www.sciencedirect.com/book/9780128029992/the-yaws-handbook-of-vapor-pressure">https://www.sciencedirect.com/book/9780128029992/the-yaws-handbook-of-vapor-pressure</a> |

## Legend

|                 |   |
|-----------------|---|
| <b>cpg:</b>     | Ideal gas heat capacity                         |
| <b>gf:</b>      | Standard Gibbs free energy of formation         |
| <b>hf:</b>      | Enthalpy of formation at standard conditions    |
| <b>hfus:</b>    | Enthalpy of fusion at standard conditions       |
| <b>hvap:</b>    | Enthalpy of vaporization at standard conditions |
| <b>log10ws:</b> | Log10 of Water solubility in mol/l              |
| <b>logp:</b>    | Octanol/Water partition coefficient             |
| <b>mcvol:</b>   | McGowan's characteristic volume                 |
| <b>pc:</b>      | Critical Pressure                               |
| <b>pvap:</b>    | Vapor pressure                                  |
| <b>rinpola:</b> | Non-polar retention indices                     |

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

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