

# Cyclohexane, 2-(1-decylundecyl)-1,4-dimethyl-

|                      |                                                                                   |
|----------------------|-----------------------------------------------------------------------------------|
| Other names:         | 11-(2,5-Dimethylcyclohexyl)heneicosane                                            |
| Inchi:               | InChI=1S/C29H58/c1-5-7-9-11-13-15-17-19-21-28(29-25-26(3)23-24-27(29)4)22-20-18-1 |
| InchiKey:            | UPYOZECJBFFKFT-UHFFFAOYSA-N                                                       |
| Formula:             | C29H58                                                                            |
| SMILES:              | CCCCCCCCCCCC(CCCCCCCCCC)C1CC(C)CCC1C                                              |
| Mol. weight [g/mol]: | 406.77                                                                            |
| CAS:                 | 55429-27-1                                                                        |

## Physical Properties

| Property code | Value   | Unit    | Source         |
|---------------|---------|---------|----------------|
| gf            | 199.89  | kJ/mol  | Joback Method  |
| hf            | -633.53 | kJ/mol  | Joback Method  |
| hfus          | 61.32   | kJ/mol  | Joback Method  |
| hvap          | 79.57   | kJ/mol  | Joback Method  |
| log10ws       | -10.89  |         | Crippen Method |
| logp          | 10.736  |         | Crippen Method |
| mcvol         | 408.610 | ml/mol  | McGowan Method |
| pc            | 672.90  | kPa     | Joback Method  |
| tb            | 872.69  | K       | Joback Method  |
| tc            | 1068.43 | K       | Joback Method  |
| tf            | 400.49  | K       | Joback Method  |
| vc            | 1.585   | m3/kmol | Joback Method  |

## Temperature Dependent Properties

| Property code | Value     | Unit    | Temperature [K] | Source        |
|---------------|-----------|---------|-----------------|---------------|
| cpg           | 1421.60   | J/mol×K | 872.69          | Joback Method |
| cpg           | 1536.45   | J/mol×K | 1035.81         | Joback Method |
| cpg           | 1516.43   | J/mol×K | 1003.18         | Joback Method |
| cpg           | 1494.99   | J/mol×K | 970.56          | Joback Method |
| cpg           | 1472.08   | J/mol×K | 937.94          | Joback Method |
| cpg           | 1447.64   | J/mol×K | 905.31          | Joback Method |
| cpg           | 1555.11   | J/mol×K | 1068.43         | Joback Method |
| dvisc         | 0.0000453 | Pa×s    | 872.69          | Joback Method |

|       |           |      |        |               |
|-------|-----------|------|--------|---------------|
| dvisc | 0.0000616 | Pa×s | 793.99 | Joback Method |
| dvisc | 0.0000896 | Pa×s | 715.29 | Joback Method |
| dvisc | 0.0001431 | Pa×s | 636.59 | Joback Method |
| dvisc | 0.0002607 | Pa×s | 557.89 | Joback Method |
| dvisc | 0.0005786 | Pa×s | 479.19 | Joback Method |
| dvisc | 0.0017563 | Pa×s | 400.49 | Joback Method |

## Sources

|                        |                                                                                                                                               |
|------------------------|-----------------------------------------------------------------------------------------------------------------------------------------------|
| <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=C55429271&amp;Units=SI">http://webbook.nist.gov/cgi/cbook.cgi?ID=C55429271&amp;Units=SI</a> |
| <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>                             |

## Legend

|                 |                                                 |
|-----------------|-------------------------------------------------|
| <b>cpg:</b>     | Ideal gas heat capacity                         |
| <b>dvisc:</b>   | Dynamic viscosity                               |
| <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>tb:</b>      | Normal Boiling Point Temperature                |
| <b>tc:</b>      | Critical Temperature                            |
| <b>tf:</b>      | Normal melting (fusion) point                   |
| <b>vc:</b>      | Critical Volume                                 |

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