

# 9-Decenylcyclopentane

|                             |   |
|-----------------------------|---|
| <b>Inchi:</b>               | InChI=1S/C15H28/c1-2-3-4-5-6-7-8-9-12-15-13-10-11-14-15/h2,15H,1,3-14H2 |
| <b>InchiKey:</b>            | HSGXSXQEZDGBHC-UHFFFAOYSA-N   |
| <b>Formula:</b>             | C15H28  |
| <b>SMILES:</b>              | C=CCCCCCCCC1CCCC1   |
| <b>Mol. weight [g/mol]:</b> | 208.38  |

## Physical Properties

| Property code | Value   | Unit                 | Source         |
|---------------|---------|----------------------|----------------|
| gf            | 199.81  | kJ/mol               | Joback Method  |
| hf            | -167.02 | kJ/mol               | Joback Method  |
| hfus          | 27.26   | kJ/mol               | Joback Method  |
| hvap          | 48.57   | kJ/mol               | Joback Method  |
| log10ws       | -5.61   |                      | Crippen Method |
| logp          | 5.483   |                      | Crippen Method |
| mcvol         | 207.050 | ml/mol               | McGowan Method |
| pc            | 1685.18 | kPa                  | Joback Method  |
| rinsol        | 1524.90 |                      | NIST Webbook   |
| tb            | 554.56  | K                    | Joback Method  |
| tc            | 737.61  | K                    | Joback Method  |
| tf            | 267.95  | K                    | Joback Method  |
| vc            | 0.797   | m <sup>3</sup> /kmol | Joback Method  |

## Temperature Dependent Properties

| Property code | Value     | Unit    | Temperature [K] | Source        |
|---------------|-----------|---------|-----------------|---------------|
| cpg           | 528.46    | J/molxK | 554.56          | Joback Method |
| cpg           | 623.09    | J/molxK | 707.10          | Joback Method |
| cpg           | 606.06    | J/molxK | 676.59          | Joback Method |
| cpg           | 588.11    | J/molxK | 646.09          | Joback Method |
| cpg           | 569.23    | J/molxK | 615.58          | Joback Method |
| cpg           | 549.35    | J/molxK | 585.07          | Joback Method |
| cpg           | 639.26    | J/molxK | 737.61          | Joback Method |
| dvisc         | 0.0002369 | Paxs    | 554.56          | Joback Method |
| dvisc         | 0.0003092 | Paxs    | 506.79          | Joback Method |

|       |           |      |        |               |
|-------|-----------|------|--------|---------------|
| dvisc | 0.0004267 | Paxs | 459.02 | Joback Method |
| dvisc | 0.0006346 | Paxs | 411.25 | Joback Method |
| dvisc | 0.0010476 | Paxs | 363.49 | Joback Method |
| dvisc | 0.0020126 | Paxs | 315.72 | Joback Method |
| dvisc | 0.0048801 | Paxs | 267.95 | Joback Method |

## Sources

|                        |   |
|------------------------|---|
| <b>NIST Webbook:</b>   | <a href="http://webbook.nist.gov/cgi/cbook.cgi?ID=R388900&amp;Units=SI">http://webbook.nist.gov/cgi/cbook.cgi?ID=R388900&amp;Units=SI</a> |
| <b>Crippen Method:</b> | <a href="http://pubs.acs.org/doi/abs/10.1021/ci9903071">http://pubs.acs.org/doi/abs/10.1021/ci9903071</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>                     |

## 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>rinpol:</b>  | Non-polar retention indices                     |
| <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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