

# Cyclopentane, 4-methylpentyl

|                      |                                                                 |
|----------------------|-----------------------------------------------------------------|
| Inchi:               | InChI=1S/C11H22/c1-10(2)6-5-9-11-7-3-4-8-11/h10-11H,3-9H2,1-2H3 |
| InchiKey:            | PDKAJPYRQGPDLU-UHFFFAOYSA-N                                     |
| Formula:             | C11H22                                                          |
| SMILES:              | CC(C)CCCC1CCCC1                                                 |
| Mol. weight [g/mol]: | 154.29                                                          |

## Physical Properties

| Property code | Value   | Unit    | Source         |
|---------------|---------|---------|----------------|
| gf            | 75.85   | kJ/mol  | Joback Method  |
| hf            | -215.17 | kJ/mol  | Joback Method  |
| hfus          | 14.66   | kJ/mol  | Joback Method  |
| hvap          | 39.95   | kJ/mol  | Joback Method  |
| log10ws       | -3.84   |         | Crippen Method |
| logp          | 4.003   |         | Crippen Method |
| mcvol         | 154.990 | ml/mol  | McGowan Method |
| pc            | 2298.11 | kPa     | Joback Method  |
| rinpol        | 1104.00 |         | NIST Webbook   |
| rinpol        | 1095.00 |         | NIST Webbook   |
| rinpol        | 1100.00 |         | NIST Webbook   |
| tb            | 465.92  | K       | Joback Method  |
| tc            | 657.55  | K       | Joback Method  |
| tf            | 209.63  | K       | Joback Method  |
| vc            | 0.587   | m3/kmol | Joback Method  |

## Temperature Dependent Properties

| Property code | Value  | Unit    | Temperature [K] | Source        |
|---------------|--------|---------|-----------------|---------------|
| cpg           | 346.47 | J/mol×K | 465.92          | Joback Method |
| cpg           | 366.19 | J/mol×K | 497.86          | Joback Method |
| cpg           | 384.96 | J/mol×K | 529.80          | Joback Method |
| cpg           | 402.80 | J/mol×K | 561.74          | Joback Method |
| cpg           | 419.74 | J/mol×K | 593.68          | Joback Method |
| cpg           | 435.83 | J/mol×K | 625.62          | Joback Method |
| cpg           | 451.08 | J/mol×K | 657.55          | Joback Method |

|       |           |      |        |               |
|-------|-----------|------|--------|---------------|
| dvisc | 0.0084385 | Pa×s | 209.63 | Joback Method |
| dvisc | 0.0029659 | Pa×s | 252.34 | Joback Method |
| dvisc | 0.0014110 | Pa×s | 295.06 | Joback Method |
| dvisc | 0.0008100 | Pa×s | 337.77 | Joback Method |
| dvisc | 0.0005267 | Pa×s | 380.49 | Joback Method |
| dvisc | 0.0003736 | Pa×s | 423.21 | Joback Method |
| dvisc | 0.0002822 | Pa×s | 465.92 | Joback Method |

## Sources

|                        |                                                                                                                                         |
|------------------------|-----------------------------------------------------------------------------------------------------------------------------------------|
| <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=R10735&amp;Units=SI">http://webbook.nist.gov/cgi/cbook.cgi?ID=R10735&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>                               |

## 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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