

# 7-Cyclohexylnonadecane

|                             |  |
|-----------------------------|--|
| <b>Inchi:</b>               | InChI=1S/C25H50/c1-3-5-7-9-10-11-12-13-14-17-21-24(20-16-8-6-4-2)25-22-18-15-19-23 |
| <b>InchiKey:</b>            | AQQBKKMWNJHQPX-UHFFFAOYSA-N  |
| <b>Formula:</b>             | C25H50   |
| <b>SMILES:</b>              | CCCCCCCCCCCC(CCCCC)C1CCCCC1  |
| <b>Mol. weight [g/mol]:</b> | 350.66   |

## Physical Properties

| Property code | Value   | Unit                 | Source         |
|---------------|---------|----------------------|----------------|
| gf            | 181.63  | kJ/mol               | Joback Method  |
| hf            | -510.29 | kJ/mol               | Joback Method  |
| hfus          | 48.82   | kJ/mol               | Joback Method  |
| hvap          | 71.28   | kJ/mol               | Joback Method  |
| log10ws       | -9.70   |                      | Crippen Method |
| logp          | 9.464   |                      | Crippen Method |
| mcvol         | 352.250 | ml/mol               | McGowan Method |
| pc            | 862.51  | kPa                  | Joback Method  |
| rinpol        | 2443.00 |                      | NIST Webbook   |
| ripol         | 2502.12 |                      | NIST Webbook   |
| tb            | 790.51  | K                    | Joback Method  |
| tc            | 974.21  | K                    | Joback Method  |
| tf            | 363.89  | K                    | Joback Method  |
| vc            | 1.363   | m <sup>3</sup> /kmol | Joback Method  |

## Temperature Dependent Properties

| Property code | Value     | Unit    | Temperature [K] | Source        |
|---------------|-----------|---------|-----------------|---------------|
| cpg           | 1152.87   | J/molxK | 790.51          | Joback Method |
| cpg           | 1261.82   | J/molxK | 943.60          | Joback Method |
| cpg           | 1242.42   | J/molxK | 912.98          | Joback Method |
| cpg           | 1221.88   | J/molxK | 882.36          | Joback Method |
| cpg           | 1200.14   | J/molxK | 851.74          | Joback Method |
| cpg           | 1177.15   | J/molxK | 821.13          | Joback Method |
| cpg           | 1280.12   | J/molxK | 974.21          | Joback Method |
| dvisc         | 0.0000482 | Paxs    | 790.51          | Joback Method |

|       |           |      |        |               |
|-------|-----------|------|--------|---------------|
| dvisc | 0.0000686 | Paxs | 719.41 | Joback Method |
| dvisc | 0.0001056 | Paxs | 648.30 | Joback Method |
| dvisc | 0.0001806 | Paxs | 577.20 | Joback Method |
| dvisc | 0.0003594 | Paxs | 506.10 | Joback Method |
| dvisc | 0.0008954 | Paxs | 434.99 | Joback Method |
| dvisc | 0.0031872 | Paxs | 363.89 | 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=U357254&amp;Units=SI">http://webbook.nist.gov/cgi/cbook.cgi?ID=U357254&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>ripol:</b>   | 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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