

# 7,11-Dimethylhexacosane

**Inchi:** InChI=1S/C28H58/c1-5-7-9-11-12-13-14-15-16-17-18-19-21-24-28(4)26-22-25-27(3)23-2  
**InchiKey:** WFTYRXXYPQHOFK-UHFFFAOYSA-N  
**Formula:** C28H58  
**SMILES:** CCCCCCCCCCCCCCCC(C)CCCC(C)CCCCC  
**Mol. weight [g/mol]:** 394.76

## Physical Properties

| Property code | Value   | Unit                 | Source         |
|---------------|---------|----------------------|----------------|
| gf            | 180.00  | kJ/mol               | Joback Method  |
| hf            | -631.81 | kJ/mol               | Joback Method  |
| hfus          | 61.23   | kJ/mol               | Joback Method  |
| hvap          | 77.15   | kJ/mol               | Joback Method  |
| log10ws       | -11.06  |                      | Crippen Method |
| logp          | 10.881  |                      | Crippen Method |
| mcvol         | 405.380 | ml/mol               | McGowan Method |
| pc            | 660.85  | kPa                  | Joback Method  |
| rinpol        | 2675.00 |                      | NIST Webbook   |
| rinpol        | 2675.00 |                      | NIST Webbook   |
| tb            | 839.16  | K                    | Joback Method  |
| tc            | 1028.00 | K                    | Joback Method  |
| tf            | 375.32  | K                    | Joback Method  |
| vc            | 1.591   | m <sup>3</sup> /kmol | Joback Method  |

## Temperature Dependent Properties

| Property code | Value     | Unit    | Temperature [K] | Source        |
|---------------|-----------|---------|-----------------|---------------|
| cpg           | 1348.71   | J/molxK | 839.16          | Joback Method |
| cpg           | 1374.05   | J/molxK | 870.63          | Joback Method |
| cpg           | 1398.06   | J/molxK | 902.11          | Joback Method |
| cpg           | 1420.81   | J/molxK | 933.58          | Joback Method |
| cpg           | 1442.36   | J/molxK | 965.06          | Joback Method |
| cpg           | 1462.77   | J/molxK | 996.53          | Joback Method |
| cpg           | 1482.09   | J/molxK | 1028.00         | Joback Method |
| dvisc         | 0.0023569 | Paxs    | 375.32          | Joback Method |

|       |           |      |        |               |
|-------|-----------|------|--------|---------------|
| dvisc | 0.0005956 | Paxs | 452.63 | Joback Method |
| dvisc | 0.0002249 | Paxs | 529.93 | Joback Method |
| dvisc | 0.0001088 | Paxs | 607.24 | Joback Method |
| dvisc | 0.0000620 | Paxs | 684.55 | Joback Method |
| dvisc | 0.0000396 | Paxs | 761.85 | Joback Method |
| dvisc | 0.0000275 | Paxs | 839.16 | Joback Method |

## Sources

|                        |   |
|------------------------|---|
| <b>NIST Webbook:</b>   | <a href="http://webbook.nist.gov/cgi/cbook.cgi?ID=R505769&amp;Units=SI">http://webbook.nist.gov/cgi/cbook.cgi?ID=R505769&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                                 |

Latest version available from:

<https://www.chemeo.com/cid/57-732-6/7-11-Dimethylhexacosane.pdf>

Generated by Cheméo on 2024-04-29 11:30:47.441748667 +0000 UTC m=+16679496.362325978.

Cheméo (<https://www.chemeo.com>) is the biggest free database of chemical and physical data for the process industry.