

# Phenethyl palmitate

|                             |   |
|-----------------------------|---|
| <b>Inchi:</b>               | InChI=1S/C24H40O2/c1-2-3-4-5-6-7-8-9-10-11-12-13-17-20-24(25)26-22-21-23-18-15-14 |
| <b>InchiKey:</b>            | PMYIGBNCUHWDMR-UHFFFAOYSA-N   |
| <b>Formula:</b>             | C24H40O2  |
| <b>SMILES:</b>              | CCCCCCCCCCCCCCCC(=O)OCCc1ccccc1   |
| <b>Mol. weight [g/mol]:</b> | 360.57  |
| <b>CAS:</b>                 | 72934-12-4  |

## Physical Properties

| Property code | Value   | Unit                 | Source         |
|---------------|---------|----------------------|----------------|
| gf            | 29.69   | kJ/mol               | Joback Method  |
| hf            | -546.96 | kJ/mol               | Joback Method  |
| hfus          | 54.74   | kJ/mol               | Joback Method  |
| hvap          | 80.45   | kJ/mol               | Joback Method  |
| log10ws       | -7.83   |                      | Crippen Method |
| logp          | 7.254   |                      | Crippen Method |
| mvol          | 332.700 | ml/mol               | McGowan Method |
| pc            | 1011.02 | kPa                  | Joback Method  |
| rinpol        | 2693.00 |                      | NIST Webbook   |
| rinpol        | 2693.00 |                      | NIST Webbook   |
| tb            | 851.49  | K                    | Joback Method  |
| tc            | 1046.47 | K                    | Joback Method  |
| tf            | 458.82  | K                    | Joback Method  |
| vc            | 1.296   | m <sup>3</sup> /kmol | Joback Method  |

## Temperature Dependent Properties

| Property code | Value   | Unit    | Temperature [K] | Source        |
|---------------|---------|---------|-----------------|---------------|
| cpg           | 1067.64 | J/molxK | 851.49          | Joback Method |
| cpg           | 1086.92 | J/molxK | 883.99          | Joback Method |
| cpg           | 1105.03 | J/molxK | 916.48          | Joback Method |
| cpg           | 1122.01 | J/molxK | 948.98          | Joback Method |
| cpg           | 1137.91 | J/molxK | 981.47          | Joback Method |
| cpg           | 1152.77 | J/molxK | 1013.97         | Joback Method |
| cpg           | 1166.66 | J/molxK | 1046.47         | Joback Method |

|       |           |      |        |               |
|-------|-----------|------|--------|---------------|
| dvisc | 0.0008470 | Paxs | 458.82 | Joback Method |
| dvisc | 0.0003773 | Paxs | 524.26 | Joback Method |
| dvisc | 0.0002011 | Paxs | 589.71 | Joback Method |
| dvisc | 0.0001216 | Paxs | 655.15 | Joback Method |
| dvisc | 0.0000805 | Paxs | 720.60 | Joback Method |
| dvisc | 0.0000571 | Paxs | 786.04 | Joback Method |
| dvisc | 0.0000427 | Paxs | 851.49 | 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=C72934124&amp;Units=SI">http://webbook.nist.gov/cgi/cbook.cgi?ID=C72934124&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>                             |

## 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>mvol:</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/86-804-4/Phenethyl-palmitate.pdf>

Generated by Cheméo on 2024-04-23 06:51:36.408666895 +0000 UTC m=+16144345.329244207.

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