

# octyl pentacosanoate

|                             |  |
|-----------------------------|--|
| <b>Inchi:</b>               | InChI=1S/C33H66O2/c1-3-5-7-9-11-12-13-14-15-16-17-18-19-20-21-22-23-24-25-26-27- |
| <b>InchiKey:</b>            | RFPRPBHMYANGJW-UHFFFAOYSA-N  |
| <b>Formula:</b>             | C33H66O2   |
| <b>SMILES:</b>              | CCCCCCCCCCCCCCCCCCCCCCCCCCCC(=O)CCCCCCCC   |
| <b>Mol. weight [g/mol]:</b> | 494.88   |

## Physical Properties

| Property code | Value   | Unit                 | Source         |
|---------------|---------|----------------------|----------------|
| gf            | -6.94   | kJ/mol               | Joback Method  |
| hf            | -969.25 | kJ/mol               | Joback Method  |
| hfus          | 84.01   | kJ/mol               | Joback Method  |
| hvap          | 98.21   | kJ/mol               | Joback Method  |
| log10ws       | -12.50  |                      | Crippen Method |
| logp          | 11.882  |                      | Crippen Method |
| mvol          | 483.270 | ml/mol               | McGowan Method |
| pc            | 530.91  | kPa                  | Joback Method  |
| rinpol        | 3461.54 |                      | NIST Webbook   |
| rinpol        | 3461.54 |                      | NIST Webbook   |
| tb            | 1030.73 | K                    | Joback Method  |
| tc            | 1306.79 | K                    | Joback Method  |
| tf            | 533.83  | K                    | Joback Method  |
| vc            | 1.907   | m <sup>3</sup> /kmol | Joback Method  |

## Temperature Dependent Properties

| Property code | Value     | Unit    | Temperature [K] | Source        |
|---------------|-----------|---------|-----------------|---------------|
| cpg           | 1750.75   | J/molxK | 1030.73         | Joback Method |
| cpg           | 1875.18   | J/molxK | 1260.78         | Joback Method |
| cpg           | 1854.77   | J/molxK | 1214.77         | Joback Method |
| cpg           | 1832.30   | J/molxK | 1168.76         | Joback Method |
| cpg           | 1807.60   | J/molxK | 1122.75         | Joback Method |
| cpg           | 1780.48   | J/molxK | 1076.74         | Joback Method |
| cpg           | 1893.74   | J/molxK | 1306.79         | Joback Method |
| dvisc         | 0.0000105 | Paxs    | 1030.73         | Joback Method |

|       |           |      |        |               |
|-------|-----------|------|--------|---------------|
| dvisc | 0.0000144 | Paxs | 947.91 | Joback Method |
| dvisc | 0.0000211 | Paxs | 865.10 | Joback Method |
| dvisc | 0.0000336 | Paxs | 782.28 | Joback Method |
| dvisc | 0.0000595 | Paxs | 699.46 | Joback Method |
| dvisc | 0.0001232 | Paxs | 616.65 | Joback Method |
| dvisc | 0.0003193 | Paxs | 533.83 | Joback Method |

## Sources

|                        |   |
|------------------------|---|
| <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=R437968&amp;Units=SI">http://webbook.nist.gov/cgi/cbook.cgi?ID=R437968&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>                                     |

## Legend

|                            |   |
|----------------------------|---|
| <b>cp<sub>g</sub>:</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>log<sub>10</sub>ws:</b> | Log <sub>10</sub> of Water solubility in mol/l  |
| <b>logp:</b>               | Octanol/Water partition coefficient             |
| <b>m<sub>cvol</sub>:</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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