

# Pimelic acid, butyl tetradecyl ester

|                             |  |
|-----------------------------|--|
| <b>Inchi:</b>               | InChI=1S/C25H48O4/c1-3-5-7-8-9-10-11-12-13-14-15-19-23-29-25(27)21-18-16-17-20-2 |
| <b>InchiKey:</b>            | MOKFGVVBNAWVMC-UHFFFAOYSA-N  |
| <b>Formula:</b>             | C25H48O4   |
| <b>SMILES:</b>              | CCCCCCCCCCCCCOC(=O)CCCCC(=O)OCCCC  |
| <b>Mol. weight [g/mol]:</b> | 412.65   |

## Physical Properties

| Property code | Value    | Unit                 | Source         |
|---------------|----------|----------------------|----------------|
| gf            | -308.22  | kJ/mol               | Joback Method  |
| hf            | -1048.93 | kJ/mol               | Joback Method  |
| hfus          | 66.08    | kJ/mol               | Joback Method  |
| hvap          | 89.56    | kJ/mol               | Joback Method  |
| log10ws       | -8.01    |                      | Crippen Method |
| logp          | 7.524    |                      | Crippen Method |
| mcvol         | 377.990  | ml/mol               | McGowan Method |
| pc            | 804.33   | kPa                  | Joback Method  |
| rinpol        | 2859.00  |                      | NIST Webbook   |
| rinpol        | 2859.00  |                      | NIST Webbook   |
| tb            | 923.98   | K                    | Joback Method  |
| tc            | 1135.00  | K                    | Joback Method  |
| tf            | 515.83   | K                    | Joback Method  |
| vc            | 1.484    | m <sup>3</sup> /kmol | Joback Method  |

## Temperature Dependent Properties

| Property code | Value     | Unit    | Temperature [K] | Source        |
|---------------|-----------|---------|-----------------|---------------|
| cpg           | 1281.17   | J/molxK | 923.98          | Joback Method |
| cpg           | 1301.94   | J/molxK | 959.15          | Joback Method |
| cpg           | 1321.17   | J/molxK | 994.32          | Joback Method |
| cpg           | 1338.92   | J/molxK | 1029.49         | Joback Method |
| cpg           | 1355.21   | J/molxK | 1064.66         | Joback Method |
| cpg           | 1370.09   | J/molxK | 1099.83         | Joback Method |
| cpg           | 1383.60   | J/molxK | 1135.00         | Joback Method |
| dvisc         | 0.0004642 | Paxs    | 515.83          | Joback Method |

|       |           |      |        |               |
|-------|-----------|------|--------|---------------|
| dvisc | 0.0002153 | Paxs | 583.86 | Joback Method |
| dvisc | 0.0001173 | Paxs | 651.88 | Joback Method |
| dvisc | 0.0000716 | Paxs | 719.90 | Joback Method |
| dvisc | 0.0000476 | Paxs | 787.93 | Joback Method |
| dvisc | 0.0000338 | Paxs | 855.95 | Joback Method |
| dvisc | 0.0000252 | Paxs | 923.98 | Joback Method |

## Sources

|                        |   |
|------------------------|---|
| <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.cheméo.com/doc/models/crippen_log10ws">https://www.cheméo.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=U406570&amp;Units=SI">http://webbook.nist.gov/cgi/cbook.cgi?ID=U406570&amp;Units=SI</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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