

# Undecanoic acid, phenylmethyl ester

|                             |   |
|-----------------------------|---|
| <b>Other names:</b>         | Benzyl undecanoate  |
| <b>Inchi:</b>               | InChI=1S/C18H28O2/c1-2-3-4-5-6-7-8-12-15-18(19)20-16-17-13-10-9-11-14-17/h9-11,13 |
| <b>InchiKey:</b>            | BBDULIAMCCXFCV-UHFFFAOYSA-N   |
| <b>Formula:</b>             | C18H28O2  |
| <b>SMILES:</b>              | CCCCCCCCCCC(=O)OCc1ccccc1   |
| <b>Mol. weight [g/mol]:</b> | 276.41  |
| <b>CAS:</b>                 | 64273-11-6  |

## Physical Properties

| Property code | Value   | Unit    | Source         |
|---------------|---------|---------|----------------|
| gf            | -20.83  | kJ/mol  | Joback Method  |
| hf            | -423.12 | kJ/mol  | Joback Method  |
| hfus          | 39.20   | kJ/mol  | Joback Method  |
| hvap          | 67.09   | kJ/mol  | Joback Method  |
| log10ws       | -5.82   |         | Crippen Method |
| logp          | 5.261   |         | Crippen Method |
| mcvol         | 248.160 | ml/mol  | McGowan Method |
| pc            | 1515.21 | kPa     | Joback Method  |
| rinpol        | 2013.00 |         | NIST Webbook   |
| tb            | 714.21  | K       | Joback Method  |
| tc            | 906.19  | K       | Joback Method  |
| tf            | 391.20  | K       | Joback Method  |
| vc            | 0.960   | m3/kmol | Joback Method  |

## Temperature Dependent Properties

| Property code | Value  | Unit    | Temperature [K] | Source        |
|---------------|--------|---------|-----------------|---------------|
| cpg           | 710.91 | J/molxK | 714.21          | Joback Method |
| cpg           | 728.69 | J/molxK | 746.21          | Joback Method |
| cpg           | 745.47 | J/molxK | 778.20          | Joback Method |
| cpg           | 761.28 | J/molxK | 810.20          | Joback Method |
| cpg           | 776.16 | J/molxK | 842.20          | Joback Method |
| cpg           | 790.14 | J/molxK | 874.20          | Joback Method |
| cpg           | 803.25 | J/molxK | 906.19          | Joback Method |

|       |           |      |        |               |
|-------|-----------|------|--------|---------------|
| dvisc | 0.0015490 | Paxs | 391.20 | Joback Method |
| dvisc | 0.0007359 | Paxs | 445.03 | Joback Method |
| dvisc | 0.0004105 | Paxs | 498.87 | Joback Method |
| dvisc | 0.0002566 | Paxs | 552.70 | Joback Method |
| dvisc | 0.0001743 | Paxs | 606.54 | Joback Method |
| dvisc | 0.0001261 | Paxs | 660.38 | Joback Method |
| dvisc | 0.0000958 | Paxs | 714.21 | 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=C64273116&amp;Units=SI">http://webbook.nist.gov/cgi/cbook.cgi?ID=C64273116&amp;Units=SI</a> |
| <b>Crippen Method:</b> | <a href="http://pubs.acs.org/doi/abs/10.1021/ci990307I">http://pubs.acs.org/doi/abs/10.1021/ci990307I</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>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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