

# Phthalic acid, butyl 8-chlorooctyl ester

|                             |  |
|-----------------------------|--|
| <b>Inchi:</b>               | InChI=1S/C20H29ClO4/c1-2-3-15-24-19(22)17-12-8-9-13-18(17)20(23)25-16-11-7-5-4-6 |
| <b>InchiKey:</b>            | RPUIJFMUHJUTCPN-UHFFFAOYSA-N   |
| <b>Formula:</b>             | C20H29ClO4   |
| <b>SMILES:</b>              | CCCCOC(=O)c1ccccc1C(=O)OCCCCCCCCCl   |
| <b>Mol. weight [g/mol]:</b> | 368.89   |

## Physical Properties

| Property code | Value   | Unit    | Source         |
|---------------|---------|---------|----------------|
| gf            | -259.47 | kJ/mol  | Joback Method  |
| hf            | -736.41 | kJ/mol  | Joback Method  |
| hfus          | 50.98   | kJ/mol  | Joback Method  |
| hvap          | 85.75   | kJ/mol  | Joback Method  |
| log10ws       | -6.30   |         | Crippen Method |
| logp          | 5.380   |         | Crippen Method |
| mcvol         | 296.020 | ml/mol  | McGowan Method |
| pc            | 1306.12 | kPa     | Joback Method  |
| rinpol        | 2678.00 |         | NIST Webbook   |
| tb            | 878.67  | K       | Joback Method  |
| tc            | 1084.02 | K       | Joback Method  |
| tf            | 528.34  | K       | Joback Method  |
| vc            | 1.145   | m3/kmol | Joback Method  |

## Temperature Dependent Properties

| Property code | Value     | Unit    | Temperature [K] | Source        |
|---------------|-----------|---------|-----------------|---------------|
| cpg           | 904.96    | J/molxK | 878.67          | Joback Method |
| cpg           | 919.93    | J/molxK | 912.90          | Joback Method |
| cpg           | 933.75    | J/molxK | 947.12          | Joback Method |
| cpg           | 946.44    | J/molxK | 981.35          | Joback Method |
| cpg           | 958.03    | J/molxK | 1015.57         | Joback Method |
| cpg           | 968.55    | J/molxK | 1049.80         | Joback Method |
| cpg           | 978.02    | J/molxK | 1084.02         | Joback Method |
| dvisc         | 0.0004842 | Paxs    | 528.34          | Joback Method |
| dvisc         | 0.0002709 | Paxs    | 586.73          | Joback Method |

|       |           |      |        |               |
|-------|-----------|------|--------|---------------|
| dvisc | 0.0001684 | Paxs | 645.12 | Joback Method |
| dvisc | 0.0001132 | Paxs | 703.50 | Joback Method |
| dvisc | 0.0000809 | Paxs | 761.89 | Joback Method |
| dvisc | 0.0000607 | Paxs | 820.28 | Joback Method |
| dvisc | 0.0000473 | Paxs | 878.67 | 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.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>                     |
| <b>NIST Webbook:</b>   | <a href="http://webbook.nist.gov/cgi/cbook.cgi?ID=U356859&amp;Units=SI">http://webbook.nist.gov/cgi/cbook.cgi?ID=U356859&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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