

# Diglycolic acid, 3,7-dimethyloctyl propyl ester

|                             |   |
|-----------------------------|---|
| <b>Inchi:</b>               | InChI=1S/C17H32O5/c1-5-10-21-16(18)12-20-13-17(19)22-11-9-15(4)8-6-7-14(2)3/h14-1 |
| <b>InchiKey:</b>            | ZVJHDPJEHMEELRC-UHFFFAOYSA-N  |
| <b>Formula:</b>             | C17H32O5  |
| <b>SMILES:</b>              | CCCOC(=O)COCC(=O)OCCC(C)CCCC(C)C  |
| <b>Mol. weight [g/mol]:</b> | 316.43  |

## Physical Properties

| Property code | Value    | Unit                 | Source         |
|---------------|----------|----------------------|----------------|
| gf            | -485.46  | kJ/mol               | Joback Method  |
| hf            | -1026.59 | kJ/mol               | Joback Method  |
| hfus          | 39.50    | kJ/mol               | Joback Method  |
| hvap          | 73.38    | kJ/mol               | Joback Method  |
| log10ws       | -3.27    |                      | Crippen Method |
| logp          | 3.352    |                      | Crippen Method |
| mcvol         | 271.140  | ml/mol               | McGowan Method |
| pc            | 1313.70  | kPa                  | Joback Method  |
| rinpol        | 2515.00  |                      | NIST Webbook   |
| rinpol        | 2515.00  |                      | NIST Webbook   |
| tb            | 762.48   | K                    | Joback Method  |
| tc            | 944.32   | K                    | Joback Method  |
| tf            | 417.90   | K                    | Joback Method  |
| vc            | 1.042    | m <sup>3</sup> /kmol | Joback Method  |

## Temperature Dependent Properties

| Property code | Value     | Unit    | Temperature [K] | Source        |
|---------------|-----------|---------|-----------------|---------------|
| cpg           | 822.56    | J/mol×K | 762.48          | Joback Method |
| cpg           | 898.47    | J/mol×K | 914.02          | Joback Method |
| cpg           | 885.19    | J/mol×K | 883.71          | Joback Method |
| cpg           | 870.96    | J/mol×K | 853.40          | Joback Method |
| cpg           | 855.78    | J/mol×K | 823.09          | Joback Method |
| cpg           | 839.65    | J/mol×K | 792.79          | Joback Method |
| cpg           | 910.80    | J/mol×K | 944.32          | Joback Method |
| dvisc         | 0.0000520 | Paxs    | 762.48          | Joback Method |

|       |           |      |        |               |
|-------|-----------|------|--------|---------------|
| dvisc | 0.0000705 | Paxs | 705.05 | Joback Method |
| dvisc | 0.0001008 | Paxs | 647.62 | Joback Method |
| dvisc | 0.0001546 | Paxs | 590.19 | Joback Method |
| dvisc | 0.0002600 | Paxs | 532.76 | Joback Method |
| dvisc | 0.0004957 | Paxs | 475.33 | Joback Method |
| dvisc | 0.0011284 | Paxs | 417.90 | 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=U382143&amp;Units=SI">http://webbook.nist.gov/cgi/cbook.cgi?ID=U382143&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>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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