

# Diglycolic acid, isobutyl oct-4-yl ester

|                             |  |
|-----------------------------|--|
| <b>Inchi:</b>               | InChI=1S/C16H30O5/c1-5-7-9-14(8-6-2)21-16(18)12-19-11-15(17)20-10-13(3)4/h13-14H |
| <b>InchiKey:</b>            | WCWHIFPMHUMXHK-UHFFFAOYSA-N  |
| <b>Formula:</b>             | C16H30O5   |
| <b>SMILES:</b>              | CCCCC(CCC)OC(=O)COCC(=O)OCC(C)C  |
| <b>Mol. weight [g/mol]:</b> | 302.41   |

## Physical Properties

| Property code | Value    | Unit                 | Source         |
|---------------|----------|----------------------|----------------|
| gf            | -493.88  | kJ/mol               | Joback Method  |
| hf            | -1005.95 | kJ/mol               | Joback Method  |
| hfus          | 36.91    | kJ/mol               | Joback Method  |
| hvap          | 71.16    | kJ/mol               | Joback Method  |
| log10ws       | -3.20    |                      | Crippen Method |
| logp          | 3.104    |                      | Crippen Method |
| mvol          | 257.050  | ml/mol               | McGowan Method |
| pc            | 1410.13  | kPa                  | Joback Method  |
| rinpol        | 2287.00  |                      | NIST Webbook   |
| rinpol        | 2287.00  |                      | NIST Webbook   |
| tb            | 739.60   | K                    | Joback Method  |
| tc            | 920.56   | K                    | Joback Method  |
| tf            | 406.63   | K                    | Joback Method  |
| vc            | 0.986    | m <sup>3</sup> /kmol | Joback Method  |

## Temperature Dependent Properties

| Property code | Value     | Unit    | Temperature [K] | Source        |
|---------------|-----------|---------|-----------------|---------------|
| cpg           | 764.62    | J/molxK | 739.60          | Joback Method |
| cpg           | 839.25    | J/molxK | 890.40          | Joback Method |
| cpg           | 826.12    | J/molxK | 860.24          | Joback Method |
| cpg           | 812.09    | J/molxK | 830.08          | Joback Method |
| cpg           | 797.17    | J/molxK | 799.92          | Joback Method |
| cpg           | 781.34    | J/molxK | 769.76          | Joback Method |
| cpg           | 851.49    | J/molxK | 920.56          | Joback Method |
| dvisc         | 0.0000599 | Paxs    | 739.60          | Joback Method |

|       |           |      |        |               |
|-------|-----------|------|--------|---------------|
| dvisc | 0.0000810 | Paxs | 684.11 | Joback Method |
| dvisc | 0.0001156 | Paxs | 628.61 | Joback Method |
| dvisc | 0.0001766 | Paxs | 573.12 | Joback Method |
| dvisc | 0.0002955 | Paxs | 517.62 | Joback Method |
| dvisc | 0.0005594 | Paxs | 462.12 | Joback Method |
| dvisc | 0.0012606 | Paxs | 406.63 | 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=U382019&amp;Units=SI">http://webbook.nist.gov/cgi/cbook.cgi?ID=U382019&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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