

# Diglycolic acid, hexyl oct-4-yl ester

|                             |   |
|-----------------------------|---|
| <b>Inchi:</b>               | InChI=1S/C18H34O5/c1-4-7-9-10-13-22-17(19)14-21-15-18(20)23-16(11-6-3)12-8-5-2/h1 |
| <b>InchiKey:</b>            | BDPSLBOHSSABGT-UHFFFAOYSA-N   |
| <b>Formula:</b>             | C18H34O5  |
| <b>SMILES:</b>              | CCCCCOC(=O)COCC(=O)OC(CCC)CCCC  |
| <b>Mol. weight [g/mol]:</b> | 330.46  |

## Physical Properties

| Property code | Value    | Unit                 | Source         |
|---------------|----------|----------------------|----------------|
| gf            | -474.60  | kJ/mol               | Joback Method  |
| hf            | -1041.95 | kJ/mol               | Joback Method  |
| hfus          | 45.61    | kJ/mol               | Joback Method  |
| hvap          | 76.00    | kJ/mol               | Joback Method  |
| log10ws       | -4.28    |                      | Crippen Method |
| logp          | 4.029    |                      | Crippen Method |
| mcvol         | 285.230  | ml/mol               | McGowan Method |
| pc            | 1219.99  | kPa                  | Joback Method  |
| rinpol        | 2579.00  |                      | NIST Webbook   |
| rinpol        | 2579.00  |                      | NIST Webbook   |
| tb            | 785.80   | K                    | Joback Method  |
| tc            | 967.95   | K                    | Joback Method  |
| tf            | 444.17   | K                    | Joback Method  |
| vc            | 1.103    | m <sup>3</sup> /kmol | Joback Method  |

## Temperature Dependent Properties

| Property code | Value     | Unit    | Temperature [K] | Source        |
|---------------|-----------|---------|-----------------|---------------|
| cpg           | 880.97    | J/molxK | 785.80          | Joback Method |
| cpg           | 898.29    | J/molxK | 816.16          | Joback Method |
| cpg           | 914.61    | J/molxK | 846.52          | Joback Method |
| cpg           | 929.94    | J/molxK | 876.87          | Joback Method |
| cpg           | 944.29    | J/molxK | 907.23          | Joback Method |
| cpg           | 957.65    | J/molxK | 937.59          | Joback Method |
| cpg           | 970.03    | J/molxK | 967.95          | Joback Method |
| dvisc         | 0.0008380 | Paxs    | 444.17          | Joback Method |

|       |           |      |        |               |
|-------|-----------|------|--------|---------------|
| dvisc | 0.0003989 | Paxs | 501.11 | Joback Method |
| dvisc | 0.0002209 | Paxs | 558.05 | Joback Method |
| dvisc | 0.0001365 | Paxs | 614.98 | Joback Method |
| dvisc | 0.0000915 | Paxs | 671.92 | Joback Method |
| dvisc | 0.0000653 | Paxs | 728.86 | Joback Method |
| dvisc | 0.0000489 | Paxs | 785.80 | 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=U382023&amp;Units=SI">http://webbook.nist.gov/cgi/cbook.cgi?ID=U382023&amp;Units=SI</a> |
| <b>Crippen Method:</b> | <a href="http://pubs.acs.org/doi/abs/10.1021/ci990307l">http://pubs.acs.org/doi/abs/10.1021/ci990307l</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>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                                 |

Latest version available from:

<https://www.chemeo.com/cid/81-844-5/Diglycolic-acid-hexyl-oct-4-yl-ester.pdf>

Generated by Cheméo on 2024-04-29 03:50:07.122538147 +0000 UTC m=+16651856.043115469.

Cheméo (<https://www.chemeo.com>) is the biggest free database of chemical and physical data for the process industry.