

# Diglycolic acid, pentyl undecyl ester

Inchi:

InchiKey:

Formula:

SMILES:

Mol. weight [g/mol]:

InChI=1S/C20H38O5/c1-3-5-7-8-9-10-11-12-14-16-25-20(22)18-23-17-19(21)24-15-13-6

HLTFLLLXQPTJFB-UHFFFAOYSA-N

C20H38O5

CCCCCCCCCCCCOC(=O)COCC(=O)OCCCCC

358.51

## Physical Properties

| Property code | Value    | Unit    | Source         |
|---------------|----------|---------|----------------|
| gf            | -455.32  | kJ/mol  | Joback Method  |
| hf            | -1077.95 | kJ/mol  | Joback Method  |
| hfus          | 54.32    | kJ/mol  | Joback Method  |
| hvap          | 80.84    | kJ/mol  | Joback Method  |
| log10ws       | -5.01    |         | Crippen Method |
| logp          | 4.810    |         | Crippen Method |
| mcvol         | 313.410  | ml/mol  | McGowan Method |
| pc            | 1065.87  | kPa     | Joback Method  |
| rinpol        | 3026.00  |         | NIST Webbook   |
| rinpol        | 3026.00  |         | NIST Webbook   |
| tb            | 832.00   | K       | Joback Method  |
| tc            | 1019.41  | K       | Joback Method  |
| tf            | 481.71   | K       | Joback Method  |
| vc            | 1.222    | m3/kmol | Joback Method  |

## Temperature Dependent Properties

| Property code | Value     | Unit    | Temperature [K] | Source        |
|---------------|-----------|---------|-----------------|---------------|
| cpg           | 1000.84   | J/mol×K | 832.00          | Joback Method |
| cpg           | 1018.88   | J/mol×K | 863.24          | Joback Method |
| cpg           | 1035.80   | J/mol×K | 894.47          | Joback Method |
| cpg           | 1051.59   | J/mol×K | 925.71          | Joback Method |
| cpg           | 1066.26   | J/mol×K | 956.94          | Joback Method |
| cpg           | 1079.84   | J/mol×K | 988.18          | Joback Method |
| cpg           | 1092.31   | J/mol×K | 1019.41         | Joback Method |
| dvisc         | 0.0005736 | Pa×s    | 481.71          | Joback Method |

|       |           |      |        |               |
|-------|-----------|------|--------|---------------|
| dvisc | 0.0002891 | Pa×s | 540.09 | Joback Method |
| dvisc | 0.0001666 | Pa×s | 598.47 | Joback Method |
| dvisc | 0.0001059 | Pa×s | 656.86 | Joback Method |
| dvisc | 0.0000725 | Pa×s | 715.24 | Joback Method |
| dvisc | 0.0000525 | Pa×s | 773.62 | Joback Method |
| dvisc | 0.0000398 | Pa×s | 832.00 | Joback Method |

## Sources

|                        |                                                                                                                                           |
|------------------------|-------------------------------------------------------------------------------------------------------------------------------------------|
| <b>NIST Webbook:</b>   | <a href="http://webbook.nist.gov/cgi/cbook.cgi?ID=U381864&amp;Units=SI">http://webbook.nist.gov/cgi/cbook.cgi?ID=U381864&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>                         |
| <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>                     |

## 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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