

# Glutaric acid, hexyl 1-phenylpropyl ester

|                             |   |
|-----------------------------|---|
| <b>Inchi:</b>               | InChI=1S/C20H30O4/c1-3-5-6-10-16-23-19(21)14-11-15-20(22)24-18(4-2)17-12-8-7-9-13 |
| <b>InchiKey:</b>            | ODEZBRVNBBDTQHQ-UHFFFAOYSA-N  |
| <b>Formula:</b>             | C20H30O4  |
| <b>SMILES:</b>              | CCCCCOC(=O)CCCC(=O)OC(CC)c1ccccc1   |
| <b>Mol. weight [g/mol]:</b> | 334.45  |

## Physical Properties

| Property code | Value   | Unit                 | Source         |
|---------------|---------|----------------------|----------------|
| gf            | -240.35 | kJ/mol               | Joback Method  |
| hf            | -714.48 | kJ/mol               | Joback Method  |
| hfus          | 43.65   | kJ/mol               | Joback Method  |
| hvap          | 80.31   | kJ/mol               | Joback Method  |
| log10ws       | -5.48   |                      | Crippen Method |
| logp          | 4.975   |                      | Crippen Method |
| mvol          | 283.780 | ml/mol               | McGowan Method |
| pc            | 1364.66 | kPa                  | Joback Method  |
| rinpol        | 2378.00 |                      | NIST Webbook   |
| tb            | 835.82  | K                    | Joback Method  |
| tc            | 1036.72 | K                    | Joback Method  |
| tf            | 470.90  | K                    | Joback Method  |
| vc            | 1.089   | m <sup>3</sup> /kmol | Joback Method  |

## Temperature Dependent Properties

| Property code | Value     | Unit    | Temperature [K] | Source        |
|---------------|-----------|---------|-----------------|---------------|
| cpg           | 878.85    | J/molxK | 835.82          | Joback Method |
| cpg           | 895.20    | J/molxK | 869.30          | Joback Method |
| cpg           | 910.39    | J/molxK | 902.79          | Joback Method |
| cpg           | 924.45    | J/molxK | 936.27          | Joback Method |
| cpg           | 937.41    | J/molxK | 969.75          | Joback Method |
| cpg           | 949.30    | J/molxK | 1003.23         | Joback Method |
| cpg           | 960.15    | J/molxK | 1036.72         | Joback Method |
| dvisc         | 0.0008026 | Paxs    | 470.90          | Joback Method |
| dvisc         | 0.0003871 | Paxs    | 531.72          | Joback Method |

|       |           |      |        |               |
|-------|-----------|------|--------|---------------|
| dvisc | 0.0002169 | Paxs | 592.54 | Joback Method |
| dvisc | 0.0001354 | Paxs | 653.36 | Joback Method |
| dvisc | 0.0000915 | Paxs | 714.18 | Joback Method |
| dvisc | 0.0000658 | Paxs | 775.00 | Joback Method |
| dvisc | 0.0000497 | Paxs | 835.82 | Joback Method |

## Sources

|                        |   |
|------------------------|---|
| <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=U358953&amp;Units=SI">http://webbook.nist.gov/cgi/cbook.cgi?ID=U358953&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>                         |
| <b>Joback Method:</b>  | <a href="https://en.wikipedia.org/wiki/Joback_method">https://en.wikipedia.org/wiki/Joback_method</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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