

# Glutaric acid, 3-ethylphenyl undecyl ester

|                             |  |
|-----------------------------|--|
| <b>Inchi:</b>               | InChI=1S/C24H38O4/c1-3-5-6-7-8-9-10-11-12-19-27-23(25)17-14-18-24(26)28-22-16-13 |
| <b>InchiKey:</b>            | HOHDNDZPKKGOGQ-UHFFFAOYSA-N  |
| <b>Formula:</b>             | C24H38O4   |
| <b>SMILES:</b>              | CCCCCCCCCOC(=O)CCCC(=O)Oc1cccc(CC)c1   |
| <b>Mol. weight [g/mol]:</b> | 390.56   |

## Physical Properties

| Property code | Value   | Unit                 | Source         |
|---------------|---------|----------------------|----------------|
| gf            | -213.86 | kJ/mol               | Joback Method  |
| hf            | -803.23 | kJ/mol               | Joback Method  |
| hfus          | 57.14   | kJ/mol               | Joback Method  |
| hvap          | 90.27   | kJ/mol               | Joback Method  |
| log10ws       | -7.31   |                      | Crippen Method |
| logp          | 6.399   |                      | Crippen Method |
| mcvol         | 340.140 | ml/mol               | McGowan Method |
| pc            | 1031.25 | kPa                  | Joback Method  |
| rinpola       | 2914.00 |                      | NIST Webbook   |
| tb            | 932.76  | K                    | Joback Method  |
| tc            | 1142.45 | K                    | Joback Method  |
| tf            | 543.50  | K                    | Joback Method  |
| vc            | 1.319   | m <sup>3</sup> /kmol | Joback Method  |

## Temperature Dependent Properties

| Property code | Value     | Unit    | Temperature [K] | Source        |
|---------------|-----------|---------|-----------------|---------------|
| cpg           | 1119.44   | J/molxK | 932.76          | Joback Method |
| cpg           | 1191.30   | J/molxK | 1107.50         | Joback Method |
| cpg           | 1179.51   | J/molxK | 1072.55         | Joback Method |
| cpg           | 1166.46   | J/molxK | 1037.60         | Joback Method |
| cpg           | 1152.13   | J/molxK | 1002.66         | Joback Method |
| cpg           | 1136.46   | J/molxK | 967.71          | Joback Method |
| cpg           | 1201.88   | J/molxK | 1142.45         | Joback Method |
| dvisc         | 0.0000308 | Paxs    | 932.76          | Joback Method |
| dvisc         | 0.0000401 | Paxs    | 867.88          | Joback Method |

|       |           |      |        |               |
|-------|-----------|------|--------|---------------|
| dvisc | 0.0000544 | Paxs | 803.01 | Joback Method |
| dvisc | 0.0000780 | Paxs | 738.13 | Joback Method |
| dvisc | 0.0001199 | Paxs | 673.25 | Joback Method |
| dvisc | 0.0002018 | Paxs | 608.38 | Joback Method |
| dvisc | 0.0003848 | Paxs | 543.50 | Joback Method |

## Sources

|                        |   |
|------------------------|---|
| <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=U359167&amp;Units=SI">http://webbook.nist.gov/cgi/cbook.cgi?ID=U359167&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>                                 |

## 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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