

# Sebacic acid, 3-methylphenyl tridecyl ester

|                             |   |
|-----------------------------|---|
| <b>Inchi:</b>               | InChI=1S/C30H50O4/c1-3-4-5-6-7-8-9-10-13-16-19-25-33-29(31)23-17-14-11-12-15-18-2 |
| <b>InchiKey:</b>            | UEPPWVODQUMDON-UHFFFAOYSA-N   |
| <b>Formula:</b>             | C30H50O4  |
| <b>SMILES:</b>              | CCCCCCCCCCCCOC(=O)CCCCCCCC(=O)Oc1cccc(C)c1  |
| <b>Mol. weight [g/mol]:</b> | 474.72  |

## Physical Properties

| Property code | Value   | Unit                 | Source         |
|---------------|---------|----------------------|----------------|
| gf            | -163.34 | kJ/mol               | Joback Method  |
| hf            | -927.07 | kJ/mol               | Joback Method  |
| hfus          | 72.68   | kJ/mol               | Joback Method  |
| hvap          | 103.62  | kJ/mol               | Joback Method  |
| log10ws       | -9.91   |                      | Crippen Method |
| logp          | 8.875   |                      | Crippen Method |
| mcvol         | 424.680 | ml/mol               | McGowan Method |
| pc            | 734.42  | kPa                  | Joback Method  |
| rinpola       | 3603.00 |                      | NIST Webbook   |
| tb            | 1070.04 | K                    | Joback Method  |
| tc            | 1324.57 | K                    | Joback Method  |
| tf            | 611.12  | K                    | Joback Method  |
| vc            | 1.655   | m <sup>3</sup> /kmol | Joback Method  |

## Temperature Dependent Properties

| Property code | Value     | Unit    | Temperature [K] | Source        |
|---------------|-----------|---------|-----------------|---------------|
| cpg           | 1496.84   | J/molxK | 1070.04         | Joback Method |
| cpg           | 1515.88   | J/molxK | 1112.46         | Joback Method |
| cpg           | 1532.83   | J/molxK | 1154.88         | Joback Method |
| cpg           | 1547.78   | J/molxK | 1197.31         | Joback Method |
| cpg           | 1560.83   | J/molxK | 1239.73         | Joback Method |
| cpg           | 1572.09   | J/molxK | 1282.15         | Joback Method |
| cpg           | 1581.64   | J/molxK | 1324.57         | Joback Method |
| dvisc         | 0.0001828 | Paxs    | 611.12          | Joback Method |
| dvisc         | 0.0000907 | Paxs    | 687.61          | Joback Method |

|       |           |      |         |               |
|-------|-----------|------|---------|---------------|
| dvisc | 0.0000517 | Paxs | 764.09  | Joback Method |
| dvisc | 0.0000327 | Paxs | 840.58  | Joback Method |
| dvisc | 0.0000223 | Paxs | 917.07  | Joback Method |
| dvisc | 0.0000161 | Paxs | 993.55  | Joback Method |
| dvisc | 0.0000122 | Paxs | 1070.04 | 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=U354935&amp;Units=SI">http://webbook.nist.gov/cgi/cbook.cgi?ID=U354935&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>                         |

## 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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