

# Sebacic acid, 2-chlorophenyl octyl ester

|                             |  |
|-----------------------------|--|
| <b>Inchi:</b>               | InChI=1S/C24H37ClO4/c1-2-3-4-5-10-15-20-28-23(26)18-11-8-6-7-9-12-19-24(27)29-22 |
| <b>InchiKey:</b>            | FXQIFOZEQNMIDW-UHFFFAOYSA-N  |
| <b>Formula:</b>             | C24H37ClO4   |
| <b>SMILES:</b>              | CCCCCCCCOC(=O)CCCCCCCC(=O)Oc1ccccc1Cl  |
| <b>Mol. weight [g/mol]:</b> | 425.00   |

## Physical Properties

| Property code | Value   | Unit                 | Source         |
|---------------|---------|----------------------|----------------|
| gf            | -225.79 | kJ/mol               | Joback Method  |
| hf            | -818.97 | kJ/mol               | Joback Method  |
| hfus          | 61.34   | kJ/mol               | Joback Method  |
| hvap          | 94.65   | kJ/mol               | Joback Method  |
| log10ws       | -8.03   |                      | Crippen Method |
| logp          | 7.270   |                      | Crippen Method |
| mvol          | 352.380 | ml/mol               | McGowan Method |
| pc            | 1007.17 | kPa                  | Joback Method  |
| rinpol        | 3144.00 |                      | NIST Webbook   |
| rinpol        | 3144.00 |                      | NIST Webbook   |
| tb            | 970.19  | K                    | Joback Method  |
| tc            | 1187.80 | K                    | Joback Method  |
| tf            | 573.42  | K                    | Joback Method  |
| vc            | 1.369   | m <sup>3</sup> /kmol | Joback Method  |

## Temperature Dependent Properties

| Property code | Value     | Unit    | Temperature [K] | Source        |
|---------------|-----------|---------|-----------------|---------------|
| cpg           | 1146.11   | J/molxK | 970.19          | Joback Method |
| cpg           | 1161.98   | J/molxK | 1006.46         | Joback Method |
| cpg           | 1176.44   | J/molxK | 1042.73         | Joback Method |
| cpg           | 1189.54   | J/molxK | 1079.00         | Joback Method |
| cpg           | 1201.32   | J/molxK | 1115.26         | Joback Method |
| cpg           | 1211.83   | J/molxK | 1151.53         | Joback Method |
| cpg           | 1221.10   | J/molxK | 1187.80         | Joback Method |
| dvisc         | 0.0003068 | Paxs    | 573.42          | Joback Method |

|       |           |      |        |               |
|-------|-----------|------|--------|---------------|
| dvisc | 0.0001644 | Paxs | 639.55 | Joback Method |
| dvisc | 0.0000990 | Paxs | 705.68 | Joback Method |
| dvisc | 0.0000650 | Paxs | 771.81 | Joback Method |
| dvisc | 0.0000457 | Paxs | 837.93 | Joback Method |
| dvisc | 0.0000338 | Paxs | 904.06 | Joback Method |
| dvisc | 0.0000260 | Paxs | 970.19 | 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=U354433&amp;Units=SI">http://webbook.nist.gov/cgi/cbook.cgi?ID=U354433&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>                                     |

## 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/95-538-0/Sebacic-acid-2-chlorophenyl-octyl-ester.pdf>

Generated by Cheméo on 2024-04-23 07:42:51.860260589 +0000 UTC m=+16147420.780837911.

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