

# Glutaric acid, heptyl 2-methyl-4-chlorophenyl ester

|                      |   |
|----------------------|---|
| Inchi:               | InChI=1S/C19H27ClO4/c1-3-4-5-6-7-13-23-18(21)9-8-10-19(22)24-17-12-11-16(20)14-15 |
| InchiKey:            | LKEQUVGCZRPIOK-UHFFFAOYSA-N   |
| Formula:             | C19H27ClO4  |
| SMILES:              | CCCCCCCOC(=O)CCCC(=O)Oc1ccc(Cl)cc1C   |
| Mol. weight [g/mol]: | 354.87  |

## Physical Properties

| Property code | Value   | Unit                 | Source         |
|---------------|---------|----------------------|----------------|
| gf            | -277.52 | kJ/mol               | Joback Method  |
| hf            | -727.24 | kJ/mol               | Joback Method  |
| hfus          | 48.00   | kJ/mol               | Joback Method  |
| hvap          | 84.18   | kJ/mol               | Joback Method  |
| log10ws       | -5.99   |                      | Crippen Method |
| logp          | 5.238   |                      | Crippen Method |
| mvol          | 281.930 | ml/mol               | McGowan Method |
| pc            | 1386.08 | kPa                  | Joback Method  |
| rinpol        | 2600.00 |                      | NIST Webbook   |
| tb            | 860.77  | K                    | Joback Method  |
| tc            | 1066.00 | K                    | Joback Method  |
| tf            | 529.59  | K                    | Joback Method  |
| vc            | 1.089   | m <sup>3</sup> /kmol | Joback Method  |

## Temperature Dependent Properties

| Property code | Value     | Unit    | Temperature [K] | Source        |
|---------------|-----------|---------|-----------------|---------------|
| cpg           | 845.41    | J/molxK | 860.77          | Joback Method |
| cpg           | 907.83    | J/molxK | 1031.80         | Joback Method |
| cpg           | 897.50    | J/molxK | 997.59          | Joback Method |
| cpg           | 886.11    | J/molxK | 963.39          | Joback Method |
| cpg           | 873.64    | J/molxK | 929.18          | Joback Method |
| cpg           | 860.08    | J/molxK | 894.98          | Joback Method |
| cpg           | 917.12    | J/molxK | 1066.00         | Joback Method |
| dvisc         | 0.0000554 | Paxs    | 860.77          | Joback Method |
| dvisc         | 0.0000699 | Paxs    | 805.57          | Joback Method |

|       |           |      |        |               |
|-------|-----------|------|--------|---------------|
| dvisc | 0.0000915 | Paxs | 750.38 | Joback Method |
| dvisc | 0.0001248 | Paxs | 695.18 | Joback Method |
| dvisc | 0.0001797 | Paxs | 639.98 | Joback Method |
| dvisc | 0.0002772 | Paxs | 584.79 | Joback Method |
| dvisc | 0.0004680 | Paxs | 529.59 | 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=U359018&amp;Units=SI">http://webbook.nist.gov/cgi/cbook.cgi?ID=U359018&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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