

# Glutaric acid, propyl 2,4,6-trichlorophenyl ester

|                      |  |
|----------------------|--|
| Inchi:               | InChI=1S/C14H15Cl3O4/c1-2-6-20-12(18)4-3-5-13(19)21-14-10(16)7-9(15)8-11(14)17/h |
| InchiKey:            | NDFXZVNVGPMJJV-UHFFFAOYSA-N  |
| Formula:             | C14H15Cl3O4  |
| SMILES:              | CCCOC(=O)CCCC(=O)Oc1c(Cl)cc(Cl)cc1Cl   |
| Mol. weight [g/mol]: | 353.62   |

## Physical Properties

| Property code | Value   | Unit    | Source         |
|---------------|---------|---------|----------------|
| gf            | -353.11 | kJ/mol  | Joback Method  |
| hf            | -666.99 | kJ/mol  | Joback Method  |
| hfus          | 43.05   | kJ/mol  | Joback Method  |
| hvap          | 82.49   | kJ/mol  | Joback Method  |
| log10ws       | -5.21   |         | Crippen Method |
| logp          | 4.676   |         | Crippen Method |
| mcvol         | 235.960 | ml/mol  | McGowan Method |
| pc            | 1898.60 | kPa     | Joback Method  |
| rinpol        | 2325.00 |         | NIST Webbook   |
| rinpol        | 2325.00 |         | NIST Webbook   |
| tb            | 826.21  | K       | Joback Method  |
| tc            | 1043.90 | K       | Joback Method  |
| tf            | 545.60  | K       | Joback Method  |
| vc            | 0.906   | m3/kmol | Joback Method  |

## Temperature Dependent Properties

| Property code | Value     | Unit    | Temperature [K] | Source        |
|---------------|-----------|---------|-----------------|---------------|
| cpg           | 609.81    | J/molxK | 826.21          | Joback Method |
| cpg           | 620.76    | J/molxK | 862.49          | Joback Method |
| cpg           | 630.78    | J/molxK | 898.77          | Joback Method |
| cpg           | 639.87    | J/molxK | 935.06          | Joback Method |
| cpg           | 648.02    | J/molxK | 971.34          | Joback Method |
| cpg           | 655.24    | J/molxK | 1007.62         | Joback Method |
| cpg           | 661.54    | J/molxK | 1043.90         | Joback Method |
| dvisc         | 0.0004707 | Paxs    | 545.60          | Joback Method |

|       |           |      |        |               |
|-------|-----------|------|--------|---------------|
| dvisc | 0.0003158 | Paxs | 592.37 | Joback Method |
| dvisc | 0.0002246 | Paxs | 639.14 | Joback Method |
| dvisc | 0.0001673 | Paxs | 685.90 | Joback Method |
| dvisc | 0.0001294 | Paxs | 732.67 | Joback Method |
| dvisc | 0.0001033 | Paxs | 779.44 | Joback Method |
| dvisc | 0.0000845 | Paxs | 826.21 | 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=U358979&amp;Units=SI">http://webbook.nist.gov/cgi/cbook.cgi?ID=U358979&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>cp<sub>g</sub>:</b>                | Ideal gas heat capacity                         |
| <b>dvisc:</b>                         | Dynamic viscosity                               |
| <b>g<sub>f</sub>:</b>                 | Standard Gibbs free energy of formation         |
| <b>h<sub>f</sub>:</b>                 | Enthalpy of formation at standard conditions    |
| <b>h<sub>fus</sub>:</b>               | Enthalpy of fusion at standard conditions       |
| <b>h<sub>vap</sub>:</b>               | Enthalpy of vaporization at standard conditions |
| <b>log<sub>10</sub>w<sub>s</sub>:</b> | Log <sub>10</sub> of Water solubility in mol/l  |
| <b>log<sub>p</sub>:</b>               | Octanol/Water partition coefficient             |
| <b>mc<sub>vol</sub>:</b>              | McGowan's characteristic volume                 |
| <b>pc:</b>                            | Critical Pressure                               |
| <b>rin<sub>pol</sub>:</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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