

# Glutaric acid, pent-4-enyl propyl ester

Inchi:

InchiKey:

Formula:

SMILES:

Mol. weight [g/mol]:

InChI=1S/C13H22O4/c1-3-5-6-11-17-13(15)9-7-8-12(14)16-10-4-2/h3H,1,4-11H2,2H3

HVTMSOCRRHZMTE-UHFFFAOYSA-N

C13H22O4

C=CCCCOC(=O)CCCC(=O)OCCC

242.31

## Physical Properties

| Property code | Value   | Unit    | Source         |
|---------------|---------|---------|----------------|
| gf            | -321.42 | kJ/mol  | Joback Method  |
| hf            | -675.82 | kJ/mol  | Joback Method  |
| hfus          | 33.72   | kJ/mol  | Joback Method  |
| hvap          | 62.17   | kJ/mol  | Joback Method  |
| log10ws       | -2.84   |         | Crippen Method |
| logp          | 2.619   |         | Crippen Method |
| mcvol         | 204.610 | ml/mol  | McGowan Method |
| pc            | 1838.84 | kPa     | Joback Method  |
| rinpol        | 1699.00 |         | NIST Webbook   |
| tb            | 646.10  | K       | Joback Method  |
| tc            | 824.98  | K       | Joback Method  |
| tf            | 378.83  | K       | Joback Method  |
| vc            | 0.792   | m3/kmol | Joback Method  |

## Temperature Dependent Properties

| Property code | Value     | Unit    | Temperature [K] | Source        |
|---------------|-----------|---------|-----------------|---------------|
| cpg           | 545.83    | J/mol×K | 646.10          | Joback Method |
| cpg           | 560.42    | J/mol×K | 675.91          | Joback Method |
| cpg           | 574.33    | J/mol×K | 705.73          | Joback Method |
| cpg           | 587.56    | J/mol×K | 735.54          | Joback Method |
| cpg           | 600.14    | J/mol×K | 765.35          | Joback Method |
| cpg           | 612.06    | J/mol×K | 795.17          | Joback Method |
| cpg           | 623.32    | J/mol×K | 824.98          | Joback Method |
| dvisc         | 0.0015761 | Pa×s    | 378.83          | Joback Method |
| dvisc         | 0.0008541 | Pa×s    | 423.38          | Joback Method |

|       |           |      |        |               |
|-------|-----------|------|--------|---------------|
| dvisc | 0.0005201 | Pa×s | 467.92 | Joback Method |
| dvisc | 0.0003453 | Pa×s | 512.46 | Joback Method |
| dvisc | 0.0002447 | Pa×s | 557.01 | Joback Method |
| dvisc | 0.0001825 | Pa×s | 601.55 | Joback Method |
| dvisc | 0.0001417 | Pa×s | 646.10 | Joback Method |

## Sources

|                        |                                                                                                                                           |
|------------------------|-------------------------------------------------------------------------------------------------------------------------------------------|
| <b>NIST Webbook:</b>   | <a href="http://webbook.nist.gov/cgi/cbook.cgi?ID=U359981&amp;Units=SI">http://webbook.nist.gov/cgi/cbook.cgi?ID=U359981&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>                                     |
| <b>McGowan Method:</b> | <a href="http://link.springer.com/article/10.1007/BF02311772">http://link.springer.com/article/10.1007/BF02311772</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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