

# Glutaric acid, butyl 3-heptyl ester

Inchi:

InchiKey:

Formula:

SMILES:

Mol. weight [g/mol]:

InChI=1S/C16H30O4/c1-4-7-10-14(6-3)20-16(18)12-9-11-15(17)19-13-8-5-2/h14H,4-13H,1-3H2

OGIFACWDKXJMLK-UHFFFAOYSA-N

C16H30O4

CCCCOC(=O)CCCC(=O)OC(CC)CCCC

286.41

## Physical Properties

| Property code | Value   | Unit    | Source         |
|---------------|---------|---------|----------------|
| gf            | -386.44 | kJ/mol  | Joback Method  |
| hf            | -868.45 | kJ/mol  | Joback Method  |
| hfus          | 39.25   | kJ/mol  | Joback Method  |
| hvap          | 69.13   | kJ/mol  | Joback Method  |
| log10ws       | -4.36   |         | Crippen Method |
| logp          | 4.012   |         | Crippen Method |
| mcvol         | 251.180 | ml/mol  | McGowan Method |
| pc            | 1419.71 | kPa     | Joback Method  |
| rinpol        | 1889.00 |         | NIST Webbook   |
| tb            | 717.62  | K       | Joback Method  |
| tc            | 895.76  | K       | Joback Method  |
| tf            | 399.40  | K       | Joback Method  |
| vc            | 0.974   | m3/kmol | Joback Method  |

## Temperature Dependent Properties

| Property code | Value     | Unit    | Temperature [K] | Source        |
|---------------|-----------|---------|-----------------|---------------|
| cpg           | 734.08    | J/mol×K | 717.62          | Joback Method |
| cpg           | 750.85    | J/mol×K | 747.31          | Joback Method |
| cpg           | 766.77    | J/mol×K | 777.00          | Joback Method |
| cpg           | 781.86    | J/mol×K | 806.69          | Joback Method |
| cpg           | 796.13    | J/mol×K | 836.38          | Joback Method |
| cpg           | 809.59    | J/mol×K | 866.07          | Joback Method |
| cpg           | 822.23    | J/mol×K | 895.76          | Joback Method |
| dvisc         | 0.0015210 | Pa×s    | 399.40          | Joback Method |
| dvisc         | 0.0007109 | Pa×s    | 452.44          | Joback Method |

|       |           |      |        |               |
|-------|-----------|------|--------|---------------|
| dvisc | 0.0003898 | Pa×s | 505.47 | Joback Method |
| dvisc | 0.0002395 | Pa×s | 558.51 | Joback Method |
| dvisc | 0.0001602 | Pa×s | 611.55 | Joback Method |
| dvisc | 0.0001142 | Pa×s | 664.58 | Joback Method |
| dvisc | 0.0000856 | Pa×s | 717.62 | 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=U359733&amp;Units=SI">http://webbook.nist.gov/cgi/cbook.cgi?ID=U359733&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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