

# Glutaric acid, hept-2-yl 3-methylbutyl ester

|                             |  |
|-----------------------------|--|
| <b>Inchi:</b>               | InChI=1S/C17H32O4/c1-5-6-7-9-15(4)21-17(19)11-8-10-16(18)20-13-12-14(2)3/h14-15H |
| <b>InchiKey:</b>            | KFHVDHKOKRZMGJ-UHFFFAOYSA-N  |
| <b>Formula:</b>             | C17H32O4   |
| <b>SMILES:</b>              | CCCCC(C)OC(=O)CCCC(=O)OCCC(C)C   |
| <b>Mol. weight [g/mol]:</b> | 300.43   |

## Physical Properties

| Property code | Value   | Unit    | Source         |
|---------------|---------|---------|----------------|
| gf            | -380.46 | kJ/mol  | Joback Method  |
| hf            | -894.37 | kJ/mol  | Joback Method  |
| hfus          | 38.31   | kJ/mol  | Joback Method  |
| hvap          | 70.97   | kJ/mol  | Joback Method  |
| log10ws       | -4.53   |         | Crippen Method |
| logp          | 4.258   |         | Crippen Method |
| mcvol         | 265.270 | ml/mol  | McGowan Method |
| pc            | 1330.04 | kPa     | Joback Method  |
| rinpol        | 1901.00 |         | NIST Webbook   |
| rinpol        | 1901.00 |         | NIST Webbook   |
| tb            | 740.06  | K       | Joback Method  |
| tc            | 920.59  | K       | Joback Method  |
| tf            | 395.67  | K       | Joback Method  |
| vc            | 1.024   | m3/kmol | Joback Method  |

## Temperature Dependent Properties

| Property code | Value     | Unit    | Temperature [K] | Source        |
|---------------|-----------|---------|-----------------|---------------|
| cpg           | 792.00    | J/molxK | 740.06          | Joback Method |
| cpg           | 869.70    | J/molxK | 890.50          | Joback Method |
| cpg           | 855.95    | J/molxK | 860.41          | Joback Method |
| cpg           | 841.31    | J/molxK | 830.32          | Joback Method |
| cpg           | 825.78    | J/molxK | 800.24          | Joback Method |
| cpg           | 809.35    | J/molxK | 770.15          | Joback Method |
| cpg           | 882.59    | J/molxK | 920.59          | Joback Method |
| dvisc         | 0.0000690 | Paxs    | 740.06          | Joback Method |

|       |           |      |        |               |
|-------|-----------|------|--------|---------------|
| dvisc | 0.0000940 | Paxs | 682.66 | Joback Method |
| dvisc | 0.0001355 | Paxs | 625.26 | Joback Method |
| dvisc | 0.0002105 | Paxs | 567.87 | Joback Method |
| dvisc | 0.0003610 | Paxs | 510.47 | Joback Method |
| dvisc | 0.0007096 | Paxs | 453.07 | Joback Method |
| dvisc | 0.0016971 | Paxs | 395.67 | 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=U392554&amp;Units=SI">http://webbook.nist.gov/cgi/cbook.cgi?ID=U392554&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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