

# Glutaric acid, nonyl 3-oxobut-2-yl ester

|                             |   |
|-----------------------------|---|
| <b>Inchi:</b>               | InChI=1S/C18H32O5/c1-4-5-6-7-8-9-10-14-22-17(20)12-11-13-18(21)23-16(3)15(2)19/h1 |
| <b>InchiKey:</b>            | OGYBFSCXBTZMEK-UHFFFAOYSA-N   |
| <b>Formula:</b>             | C18H32O5  |
| <b>SMILES:</b>              | CCCCCCCCCOC(=O)CCCC(=O)OC(C)C(C)=O  |
| <b>Mol. weight [g/mol]:</b> | 328.44  |

## Physical Properties

| Property code | Value    | Unit                 | Source         |
|---------------|----------|----------------------|----------------|
| gf            | -498.52  | kJ/mol               | Joback Method  |
| hf            | -1022.31 | kJ/mol               | Joback Method  |
| hfus          | 46.03    | kJ/mol               | Joback Method  |
| hvap          | 80.33    | kJ/mol               | Joback Method  |
| log10ws       | -4.47    |                      | Crippen Method |
| logp          | 3.971    |                      | Crippen Method |
| mcvol         | 280.930  | ml/mol               | McGowan Method |
| pc            | 1291.14  | kPa                  | Joback Method  |
| rinqol        | 2273.00  |                      | NIST Webbook   |
| tb            | 817.25   | K                    | Joback Method  |
| tc            | 1006.29  | K                    | Joback Method  |
| tf            | 471.87   | K                    | Joback Method  |
| vc            | 1.091    | m <sup>3</sup> /kmol | Joback Method  |

## Temperature Dependent Properties

| Property code | Value     | Unit    | Temperature [K] | Source        |
|---------------|-----------|---------|-----------------|---------------|
| cpg           | 872.85    | J/molxK | 817.25          | Joback Method |
| cpg           | 889.00    | J/molxK | 848.76          | Joback Method |
| cpg           | 904.13    | J/molxK | 880.26          | Joback Method |
| cpg           | 918.25    | J/molxK | 911.77          | Joback Method |
| cpg           | 931.39    | J/molxK | 943.28          | Joback Method |
| cpg           | 943.54    | J/molxK | 974.78          | Joback Method |
| cpg           | 954.72    | J/molxK | 1006.29         | Joback Method |
| dvisc         | 0.0009340 | Paxs    | 471.87          | Joback Method |
| dvisc         | 0.0004651 | Paxs    | 529.43          | Joback Method |

|       |           |      |        |               |
|-------|-----------|------|--------|---------------|
| dvisc | 0.0002655 | Paxs | 587.00 | Joback Method |
| dvisc | 0.0001676 | Paxs | 644.56 | Joback Method |
| dvisc | 0.0001140 | Paxs | 702.12 | Joback Method |
| dvisc | 0.0000823 | Paxs | 759.69 | Joback Method |
| dvisc | 0.0000621 | Paxs | 817.25 | Joback Method |

## Sources

|                        |   |
|------------------------|---|
| <b>NIST Webbook:</b>   | <a href="http://webbook.nist.gov/cgi/cbook.cgi?ID=U359709&amp;Units=SI">http://webbook.nist.gov/cgi/cbook.cgi?ID=U359709&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>                                     |
| <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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