

# Glutaric acid, dodecyl 1-naphthyl ester

**Inchi:** InChI=1S/C27H38O4/c1-2-3-4-5-6-7-8-9-10-13-22-30-26(28)20-15-21-27(29)31-25-19-14  
**InchiKey:** FYDOSLGS GPMAIL-UHFFFAOYSA-N  
**Formula:** C27H38O4  
**SMILES:** CCCCCCCCCCOC(=O)CCCC(=O)Oc1cccc2ccccc12  
**Mol. weight [g/mol]:** 426.59

## Physical Properties

| Property code | Value   | Unit                 | Source         |
|---------------|---------|----------------------|----------------|
| gf            | -81.95  | kJ/mol               | Joback Method  |
| hf            | -674.08 | kJ/mol               | Joback Method  |
| hfus          | 61.93   | kJ/mol               | Joback Method  |
| hvap          | 98.59   | kJ/mol               | Joback Method  |
| log10ws       | -8.73   |                      | Crippen Method |
| logp          | 7.380   |                      | Crippen Method |
| mvol          | 362.950 | ml/mol               | McGowan Method |
| pc            | 1003.35 | kPa                  | Joback Method  |
| rinpol        | 3449.00 |                      | NIST Webbook   |
| tb            | 1020.38 | K                    | Joback Method  |
| tc            | 1249.29 | K                    | Joback Method  |
| tf            | 610.01  | K                    | Joback Method  |
| vc            | 1.409   | m <sup>3</sup> /kmol | Joback Method  |

## Temperature Dependent Properties

| Property code | Value     | Unit    | Temperature [K] | Source        |
|---------------|-----------|---------|-----------------|---------------|
| cpg           | 1224.22   | J/molxK | 1020.38         | Joback Method |
| cpg           | 1240.48   | J/molxK | 1058.53         | Joback Method |
| cpg           | 1255.43   | J/molxK | 1096.68         | Joback Method |
| cpg           | 1269.16   | J/molxK | 1134.83         | Joback Method |
| cpg           | 1281.77   | J/molxK | 1172.99         | Joback Method |
| cpg           | 1293.34   | J/molxK | 1211.14         | Joback Method |
| cpg           | 1303.96   | J/molxK | 1249.29         | Joback Method |
| dvisc         | 0.0003583 | Paxs    | 610.01          | Joback Method |
| dvisc         | 0.0002086 | Paxs    | 678.40          | Joback Method |

|       |           |      |         |               |
|-------|-----------|------|---------|---------------|
| dvisc | 0.0001341 | Paxs | 746.80  | Joback Method |
| dvisc | 0.0000928 | Paxs | 815.19  | Joback Method |
| dvisc | 0.0000680 | Paxs | 883.59  | Joback Method |
| dvisc | 0.0000521 | Paxs | 951.98  | Joback Method |
| dvisc | 0.0000414 | Paxs | 1020.38 | 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=U358774&amp;Units=SI">http://webbook.nist.gov/cgi/cbook.cgi?ID=U358774&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>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                                 |

Latest version available from:

<https://www.chemeo.com/cid/13-756-8/Glutaric-acid-dodecyl-1-naphthyl-ester.pdf>

Generated by Cheméo on 2024-04-26 05:00:41.537283997 +0000 UTC m=+16396890.457861309.

Cheméo (<https://www.chemeo.com>) is the biggest free database of chemical and physical data for the process industry.