

# Glutaric acid, 2-ethylhexyl dodec-9-yn-1-yl ester

**Inchi:** InChI=1S/C25H44O4/c1-4-7-9-10-11-12-13-14-15-16-21-28-24(26)19-17-20-25(27)29-22  
**InchiKey:** NBAYOKJFZHLSLM-UHFFFAOYSA-N  
**Formula:** C25H44O4  
**SMILES:** CCC#CCCCCCCCCOC(=O)CCCC(=O)OCC(CC)CCCC  
**Mol. weight [g/mol]:** 408.61

## Physical Properties

| Property code | Value   | Unit    | Source         |
|---------------|---------|---------|----------------|
| gf            | -107.86 | kJ/mol  | Joback Method  |
| hf            | -781.91 | kJ/mol  | Joback Method  |
| hfus          | 65.68   | kJ/mol  | Joback Method  |
| hvap          | 91.32   | kJ/mol  | Joback Method  |
| log10ws       | -7.57   |         | Crippen Method |
| logp          | 6.604   |         | Crippen Method |
| mcvol         | 369.390 | ml/mol  | McGowan Method |
| pc            | 886.83  | kPa     | Joback Method  |
| rinpol        | 2824.00 |         | NIST Webbook   |
| rinpol        | 2824.00 |         | NIST Webbook   |
| tb            | 932.54  | K       | Joback Method  |
| tc            | 1141.86 | K       | Joback Method  |
| tf            | 606.93  | K       | Joback Method  |
| vc            | 1.440   | m3/kmol | Joback Method  |

## Temperature Dependent Properties

| Property code | Value   | Unit    | Temperature [K] | Source        |
|---------------|---------|---------|-----------------|---------------|
| cpg           | 1227.96 | J/mol×K | 932.54          | Joback Method |
| cpg           | 1247.04 | J/mol×K | 967.43          | Joback Method |
| cpg           | 1264.68 | J/mol×K | 1002.31         | Joback Method |
| cpg           | 1280.93 | J/mol×K | 1037.20         | Joback Method |
| cpg           | 1295.80 | J/mol×K | 1072.09         | Joback Method |
| cpg           | 1309.35 | J/mol×K | 1106.98         | Joback Method |
| cpg           | 1321.59 | J/mol×K | 1141.86         | Joback Method |

# Sources

|                        |                                                                                                                                           |
|------------------------|-------------------------------------------------------------------------------------------------------------------------------------------|
| <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>                     |
| <b>NIST Webbook:</b>   | <a href="http://webbook.nist.gov/cgi/cbook.cgi?ID=U393944&amp;Units=SI">http://webbook.nist.gov/cgi/cbook.cgi?ID=U393944&amp;Units=SI</a> |

# Legend

|                 |                                                 |
|-----------------|-------------------------------------------------|
| <b>cpg:</b>     | Ideal gas heat capacity                         |
| <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>rlnpol:</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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