

# Glutaric acid, 3-(2-methoxyethyl)nonyl nonyl ester

|                      |                                                                                   |
|----------------------|-----------------------------------------------------------------------------------|
| Inchi:               | InChI=1S/C26H50O5/c1-4-6-8-10-11-12-14-21-30-25(27)17-15-18-26(28)31-23-20-24(19) |
| InchiKey:            | JOACISDMSSJZJU-UHFFFAOYSA-N                                                       |
| Formula:             | C26H50O5                                                                          |
| SMILES:              | CCCCCCCCCOC(=O)CCCC(=O)OCCC(CCCCCC)CCOC                                           |
| Mol. weight [g/mol]: | 442.67                                                                            |

## Physical Properties

| Property code | Value    | Unit    | Source         |
|---------------|----------|---------|----------------|
| gf            | -407.24  | kJ/mol  | Joback Method  |
| hf            | -1207.07 | kJ/mol  | Joback Method  |
| hfus          | 66.33    | kJ/mol  | Joback Method  |
| hvap          | 93.80    | kJ/mol  | Joback Method  |
| log10ws       | -7.28    |         | Crippen Method |
| logp          | 7.007    |         | Crippen Method |
| mcvol         | 397.950  | ml/mol  | McGowan Method |
| pc            | 758.49   | kPa     | Joback Method  |
| rinpol        | 3015.00  |         | NIST Webbook   |
| rinpol        | 3015.00  |         | NIST Webbook   |
| tb            | 968.84   | K       | Joback Method  |
| tc            | 1195.64  | K       | Joback Method  |
| tf            | 534.33   | K       | Joback Method  |
| vc            | 1.552    | m3/kmol | Joback Method  |

## Temperature Dependent Properties

| Property code | Value     | Unit    | Temperature [K] | Source        |
|---------------|-----------|---------|-----------------|---------------|
| cpg           | 1377.50   | J/mol×K | 968.84          | Joback Method |
| cpg           | 1398.28   | J/mol×K | 1006.64         | Joback Method |
| cpg           | 1417.13   | J/mol×K | 1044.44         | Joback Method |
| cpg           | 1434.09   | J/mol×K | 1082.24         | Joback Method |
| cpg           | 1449.19   | J/mol×K | 1120.04         | Joback Method |
| cpg           | 1462.47   | J/mol×K | 1157.84         | Joback Method |
| cpg           | 1473.97   | J/mol×K | 1195.64         | Joback Method |
| dvisc         | 0.0003107 | Pa×s    | 534.33          | Joback Method |

|       |           |      |        |               |
|-------|-----------|------|--------|---------------|
| dvisc | 0.0001373 | Pa×s | 606.75 | Joback Method |
| dvisc | 0.0000722 | Pa×s | 679.17 | Joback Method |
| dvisc | 0.0000430 | Pa×s | 751.58 | Joback Method |
| dvisc | 0.0000281 | Pa×s | 824.00 | Joback Method |
| dvisc | 0.0000196 | Pa×s | 896.42 | Joback Method |
| dvisc | 0.0000145 | Pa×s | 968.84 | 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.cheméo.com/doc/models/crippen_log10ws">https://www.cheméo.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=U358463&amp;Units=SI">http://webbook.nist.gov/cgi/cbook.cgi?ID=U358463&amp;Units=SI</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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