

Glutaric acid, 3,3-dimethylbut-2-yl octyl ester

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|-----------------------------|---|
| Inchi: | InChI=1S/C19H36O4/c1-6-7-8-9-10-11-15-22-17(20)13-12-14-18(21)23-16(2)19(3,4)5/h1 |
| InchiKey: | QJLMHVUHXXFAHR-UHFFFAOYSA-N |
| Formula: | C19H36O4 |
| SMILES: | CCCCCCCCOC(=O)CCCC(=O)OC(C)C(C)(C)C |
| Mol. weight [g/mol]: | 328.49 |

Physical Properties

| Property code | Value | Unit | Source |
|---------------|---------|----------------------|----------------|
| gf | -358.34 | kJ/mol | Joback Method |
| hf | -939.12 | kJ/mol | Joback Method |
| hfus | 39.60 | kJ/mol | Joback Method |
| hvap | 74.52 | kJ/mol | Joback Method |
| log10ws | -5.37 | | Crippen Method |
| logp | 5.038 | | Crippen Method |
| mcvol | 293.450 | ml/mol | McGowan Method |
| pc | 1170.42 | kPa | Joback Method |
| rinqol | 2158.00 | | NIST Webbook |
| tb | 783.03 | K | Joback Method |
| tc | 968.31 | K | Joback Method |
| tf | 435.63 | K | Joback Method |
| vc | 1.131 | m ³ /kmol | Joback Method |

Temperature Dependent Properties

| Property code | Value | Unit | Temperature [K] | Source |
|---------------|-----------|---------|-----------------|---------------|
| cpg | 910.66 | J/molxK | 783.03 | Joback Method |
| cpg | 990.82 | J/molxK | 937.43 | Joback Method |
| cpg | 976.73 | J/molxK | 906.55 | Joback Method |
| cpg | 961.70 | J/molxK | 875.67 | Joback Method |
| cpg | 945.69 | J/molxK | 844.79 | Joback Method |
| cpg | 928.69 | J/molxK | 813.91 | Joback Method |
| cpg | 1004.00 | J/molxK | 968.31 | Joback Method |
| dvisc | 0.0000457 | Paxs | 783.03 | Joback Method |
| dvisc | 0.0000628 | Paxs | 725.13 | Joback Method |

| | | | | |
|-------|-----------|------|--------|---------------|
| dvisc | 0.0000912 | Paxs | 667.23 | Joback Method |
| dvisc | 0.0001421 | Paxs | 609.33 | Joback Method |
| dvisc | 0.0002430 | Paxs | 551.43 | Joback Method |
| dvisc | 0.0004713 | Paxs | 493.53 | Joback Method |
| dvisc | 0.0010903 | Paxs | 435.63 | Joback Method |

Sources

| | |
|------------------------|---|
| Crippen Method: | https://www.chemeo.com/doc/models/crippen_log10ws |
| Joback Method: | https://en.wikipedia.org/wiki/Joback_method |
| McGowan Method: | http://link.springer.com/article/10.1007/BF02311772 |
| NIST Webbook: | http://webbook.nist.gov/cgi/cbook.cgi?ID=U359755&Units=SI |
| Crippen Method: | http://pubs.acs.org/doi/abs/10.1021/ci990307l |

Legend

| | |
|-----------------|---|
| cpg: | Ideal gas heat capacity |
| dvisc: | Dynamic viscosity |
| gf: | Standard Gibbs free energy of formation |
| hf: | Enthalpy of formation at standard conditions |
| hfus: | Enthalpy of fusion at standard conditions |
| hvap: | Enthalpy of vaporization at standard conditions |
| log10ws: | Log10 of Water solubility in mol/l |
| logp: | Octanol/Water partition coefficient |
| mcvol: | McGowan's characteristic volume |
| pc: | Critical Pressure |
| rinpol: | Non-polar retention indices |
| tb: | Normal Boiling Point Temperature |
| tc: | Critical Temperature |
| tf: | Normal melting (fusion) point |
| vc: | Critical Volume |

Latest version available from:

<https://www.chemeo.com/cid/60-318-2/Glutaric-acid-3-3-dimethylbut-2-yl-octyl-ester.pdf>

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