

Glutaric acid, pent-4-enyl tetradecyl ester

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| Inchi: | InChI=1S/C24H44O4/c1-3-5-7-8-9-10-11-12-13-14-15-17-22-28-24(26)20-18-19-23(25)2 |
| InchiKey: | AGZJLYSYVVCRNT-UHFFFAOYSA-N |
| Formula: | C24H44O4 |
| SMILES: | C=CCCCOC(=O)CCCC(=O)OCCCCCCCCCCCCCCC |
| Mol. weight [g/mol]: | 396.60 |

Physical Properties

| Property code | Value | Unit | Source |
|---------------|---------|----------------------|----------------|
| gf | -228.80 | kJ/mol | Joback Method |
| hf | -902.86 | kJ/mol | Joback Method |
| hfus | 62.21 | kJ/mol | Joback Method |
| hvap | 86.66 | kJ/mol | Joback Method |
| log10ws | -7.45 | | Crippen Method |
| logp | 6.910 | | Crippen Method |
| mcvol | 359.600 | ml/mol | McGowan Method |
| pc | 871.19 | kPa | Joback Method |
| rinsol | 2804.00 | | NIST Webbook |
| tb | 897.78 | K | Joback Method |
| tc | 1099.94 | K | Joback Method |
| tf | 502.80 | K | Joback Method |
| vc | 1.409 | m ³ /kmol | Joback Method |

Temperature Dependent Properties

| Property code | Value | Unit | Temperature [K] | Source |
|---------------|-----------|---------|-----------------|---------------|
| cpg | 1188.70 | J/molxK | 897.78 | Joback Method |
| cpg | 1208.28 | J/molxK | 931.47 | Joback Method |
| cpg | 1226.53 | J/molxK | 965.17 | Joback Method |
| cpg | 1243.49 | J/molxK | 998.86 | Joback Method |
| cpg | 1259.19 | J/molxK | 1032.56 | Joback Method |
| cpg | 1273.67 | J/molxK | 1066.25 | Joback Method |
| cpg | 1286.97 | J/molxK | 1099.94 | Joback Method |
| dvisc | 0.0005395 | Paxs | 502.80 | Joback Method |
| dvisc | 0.0002558 | Paxs | 568.63 | Joback Method |

| | | | | |
|-------|-----------|------|--------|---------------|
| dvisc | 0.0001416 | Paxs | 634.46 | Joback Method |
| dvisc | 0.0000876 | Paxs | 700.29 | Joback Method |
| dvisc | 0.0000588 | Paxs | 766.12 | Joback Method |
| dvisc | 0.0000421 | Paxs | 831.95 | Joback Method |
| dvisc | 0.0000316 | Paxs | 897.78 | Joback Method |

Sources

| | |
|------------------------|-------------------------------------------------------------------------------------------------------------------------------------------|
| McGowan Method: | http://link.springer.com/article/10.1007/BF02311772 |
| NIST Webbook: | http://webbook.nist.gov/cgi/cbook.cgi?ID=U359993&Units=SI |
| Crippen Method: | http://pubs.acs.org/doi/abs/10.1021/ci9903071 |
| Crippen Method: | https://www.chemeo.com/doc/models/crippen_log10ws |
| Joback Method: | https://en.wikipedia.org/wiki/Joback_method |

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 |

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