

Glutraic acid, cis-non-3-enyl tridecyl ester

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|-----------------------------|--|
| Inchi: | InChI=1S/C27H50O4/c1-3-5-7-9-11-12-13-14-16-18-20-25-31-27(29)23-21-22-26(28)30- |
| InchiKey: | HAPBCMCEHWLBAP-ICFOKQHNSA-N |
| Formula: | C27H50O4 |
| SMILES: | CCCCC=CCCOC(=O)CCCC(=O)OCCCCCCCCCCCCC |
| Mol. weight [g/mol]: | 438.68 |

Physical Properties

| Property code | Value | Unit | Source |
|---------------|---------|---------|----------------|
| gf | -211.16 | kJ/mol | Joback Method |
| hf | -972.99 | kJ/mol | Joback Method |
| hfus | 71.46 | kJ/mol | Joback Method |
| hvap | 93.97 | kJ/mol | Joback Method |
| log10ws | -8.70 | | Crippen Method |
| logp | 8.081 | | Crippen Method |
| mcvol | 401.870 | ml/mol | McGowan Method |
| pc | 744.07 | kPa | Joback Method |
| rinpol | 3112.00 | | NIST Webbook |
| tb | 973.90 | K | Joback Method |
| tc | 1201.27 | K | Joback Method |
| tf | 533.29 | K | Joback Method |
| vc | 1.575 | m3/kmol | Joback Method |

Temperature Dependent Properties

| Property code | Value | Unit | Temperature [K] | Source |
|---------------|-----------|---------|-----------------|---------------|
| cpg | 1381.80 | J/molxK | 973.90 | Joback Method |
| cpg | 1473.28 | J/molxK | 1163.38 | Joback Method |
| cpg | 1457.97 | J/molxK | 1125.48 | Joback Method |
| cpg | 1441.24 | J/molxK | 1087.59 | Joback Method |
| cpg | 1423.02 | J/molxK | 1049.69 | Joback Method |
| cpg | 1403.23 | J/molxK | 1011.80 | Joback Method |
| cpg | 1487.25 | J/molxK | 1201.27 | Joback Method |
| dvisc | 0.0000162 | Paxs | 973.90 | Joback Method |
| dvisc | 0.0000218 | Paxs | 900.46 | Joback Method |

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|-------|-----------|------|--------|---------------|
| dvisc | 0.0000310 | Paxs | 827.03 | Joback Method |
| dvisc | 0.0000473 | Paxs | 753.60 | Joback Method |
| dvisc | 0.0000788 | Paxs | 680.16 | Joback Method |
| dvisc | 0.0001486 | Paxs | 606.72 | Joback Method |
| dvisc | 0.0003337 | Paxs | 533.29 | Joback Method |

Sources

| | |
|------------------------|---|
| NIST Webbook: | http://webbook.nist.gov/cgi/cbook.cgi?ID=U359917&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 |
| McGowan Method: | http://link.springer.com/article/10.1007/BF02311772 |

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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