

Glutaric acid, 2-octyl tridecyl ester

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| Inchi: | InChI=1S/C26H50O4/c1-4-6-8-10-11-12-13-14-15-16-18-23-29-25(27)21-19-22-26(28)30 |
| InchiKey: | BUXAUXFEUCGEAL-UHFFFAOYSA-N |
| Formula: | C26H50O4 |
| SMILES: | CCCCCCCCCCCCOC(=O)CCCC(=O)OC(C)CCCCC |
| Mol. weight [g/mol]: | 426.67 |

Physical Properties

| Property code | Value | Unit | Source |
|---------------|----------|----------------------|----------------|
| gf | -302.24 | kJ/mol | Joback Method |
| hf | -1074.85 | kJ/mol | Joback Method |
| hfus | 65.15 | kJ/mol | Joback Method |
| hvap | 91.39 | kJ/mol | Joback Method |
| log10ws | -8.54 | | Crippen Method |
| logp | 7.913 | | Crippen Method |
| mvol | 392.080 | ml/mol | McGowan Method |
| pc | 765.64 | kPa | Joback Method |
| rinpol | 2878.00 | | NIST Webbook |
| rinpol | 2878.00 | | NIST Webbook |
| tb | 946.42 | K | Joback Method |
| tc | 1164.49 | K | Joback Method |
| tf | 512.10 | K | Joback Method |
| vc | 1.534 | m ³ /kmol | Joback Method |

Temperature Dependent Properties

| Property code | Value | Unit | Temperature [K] | Source |
|---------------|-----------|---------|-----------------|---------------|
| cpg | 1345.59 | J/molxK | 946.42 | Joback Method |
| cpg | 1366.80 | J/molxK | 982.76 | Joback Method |
| cpg | 1386.34 | J/molxK | 1019.11 | Joback Method |
| cpg | 1404.25 | J/molxK | 1055.45 | Joback Method |
| cpg | 1420.58 | J/molxK | 1091.80 | Joback Method |
| cpg | 1435.40 | J/molxK | 1128.14 | Joback Method |
| cpg | 1448.73 | J/molxK | 1164.49 | Joback Method |
| dvisc | 0.0004671 | Paxs | 512.10 | Joback Method |

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|-------|-----------|------|--------|---------------|
| dvisc | 0.0001988 | Paxs | 584.49 | Joback Method |
| dvisc | 0.0001021 | Paxs | 656.87 | Joback Method |
| dvisc | 0.0000599 | Paxs | 729.26 | Joback Method |
| dvisc | 0.0000387 | Paxs | 801.65 | Joback Method |
| dvisc | 0.0000268 | Paxs | 874.03 | Joback Method |
| dvisc | 0.0000197 | Paxs | 946.42 | Joback Method |

Sources

| | |
|------------------------|---|
| 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=U358167&Units=SI |
| Crippen Method: | http://pubs.acs.org/doi/abs/10.1021/ci990307l |
| Crippen Method: | https://www.chemeo.com/doc/models/crippen_log10ws |

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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