

Sebacic acid, 2-(2-methoxyethyl)heptyl octyl ester

| | |
|----------------------|--|
| Inchi: | InChI=1S/C28H54O5/c1-4-6-8-9-14-18-23-32-27(29)20-16-12-10-11-13-17-21-28(30)33- |
| InchiKey: | VABSLKCCSHBLRL-UHFFFAOYSA-N |
| Formula: | C28H54O5 |
| SMILES: | CCCCCCCCOC(=O)CCCCCCCC(=O)OCC(CCCCC)CCOC |
| Mol. weight [g/mol]: | 470.73 |

Physical Properties

| Property code | Value | Unit | Source |
|---------------|----------|----------------------|----------------|
| gf | -390.40 | kJ/mol | Joback Method |
| hf | -1248.35 | kJ/mol | Joback Method |
| hfus | 71.52 | kJ/mol | Joback Method |
| hvap | 98.26 | kJ/mol | Joback Method |
| log10ws | -8.11 | | Crippen Method |
| logp | 7.787 | | Crippen Method |
| mvol | 426.130 | ml/mol | McGowan Method |
| pc | 684.21 | kPa | Joback Method |
| rinpol | 3216.00 | | NIST Webbook |
| tb | 1014.60 | K | Joback Method |
| tc | 1263.85 | K | Joback Method |
| tf | 556.87 | K | Joback Method |
| vc | 1.663 | m ³ /kmol | Joback Method |

Temperature Dependent Properties

| Property code | Value | Unit | Temperature [K] | Source |
|---------------|-----------|---------|-----------------|---------------|
| cpg | 1506.22 | J/molxK | 1014.60 | Joback Method |
| cpg | 1528.11 | J/molxK | 1056.14 | Joback Method |
| cpg | 1547.60 | J/molxK | 1097.68 | Joback Method |
| cpg | 1564.77 | J/molxK | 1139.22 | Joback Method |
| cpg | 1579.67 | J/molxK | 1180.76 | Joback Method |
| cpg | 1592.37 | J/molxK | 1222.30 | Joback Method |
| cpg | 1602.91 | J/molxK | 1263.85 | Joback Method |
| dvisc | 0.0002359 | Paxs | 556.87 | Joback Method |
| dvisc | 0.0001027 | Paxs | 633.16 | Joback Method |

| | | | | |
|-------|-----------|------|---------|---------------|
| dvisc | 0.0000535 | Paxs | 709.45 | Joback Method |
| dvisc | 0.0000316 | Paxs | 785.73 | Joback Method |
| dvisc | 0.0000205 | Paxs | 862.02 | Joback Method |
| dvisc | 0.0000143 | Paxs | 938.31 | Joback Method |
| dvisc | 0.0000105 | Paxs | 1014.60 | Joback Method |

Sources

| | |
|------------------------|---|
| NIST Webbook: | http://webbook.nist.gov/cgi/cbook.cgi?ID=U354357&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 |

Latest version available from:

<https://www.chemeo.com/cid/55-011-8/Sebacic-acid-2-2-methoxyethyl-heptyl-octyl-ester.pdf>

Generated by Cheméo on 2024-04-23 16:17:27.886493039 +0000 UTC m=+16178296.807070351.

Cheméo (<https://www.chemeo.com>) is the biggest free database of chemical and physical data for the process industry.