

Sebacic acid, but-2-enyl decyl ester

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|-----------------------------|--|
| Inchi: | InChI=1S/C24H44O4/c1-3-5-7-8-9-12-15-18-22-28-24(26)20-17-14-11-10-13-16-19-23(2 |
| InchiKey: | WRXAKYIPZIGLTJ-GQCTYLIASA-N |
| Formula: | C24H44O4 |
| SMILES: | CC=CCOC(=O)CCCCCCCCC(=O)OCCCCCCCCC |
| Mol. weight [g/mol]: | 396.60 |

Physical Properties

| Property code | Value | Unit | Source |
|---------------|---------|----------------------|----------------|
| gf | -236.42 | kJ/mol | Joback Method |
| hf | -911.07 | kJ/mol | Joback Method |
| hfus | 63.69 | kJ/mol | Joback Method |
| hvap | 87.29 | kJ/mol | Joback Method |
| log10ws | -7.45 | | Crippen Method |
| logp | 6.910 | | Crippen Method |
| mcvol | 359.600 | ml/mol | McGowan Method |
| pc | 876.36 | kPa | Joback Method |
| rinsol | 2772.00 | | NIST Webbook |
| tb | 905.26 | K | Joback Method |
| tc | 1108.91 | K | Joback Method |
| tf | 499.48 | K | Joback Method |
| vc | 1.407 | m ³ /kmol | Joback Method |

Temperature Dependent Properties

| Property code | Value | Unit | Temperature [K] | Source |
|---------------|-----------|---------|-----------------|---------------|
| cpg | 1190.55 | J/molxK | 905.26 | Joback Method |
| cpg | 1275.94 | J/molxK | 1074.97 | Joback Method |
| cpg | 1261.26 | J/molxK | 1041.03 | Joback Method |
| cpg | 1245.42 | J/molxK | 1007.08 | Joback Method |
| cpg | 1228.39 | J/molxK | 973.14 | Joback Method |
| cpg | 1210.11 | J/molxK | 939.20 | Joback Method |
| cpg | 1289.52 | J/molxK | 1108.91 | Joback Method |
| dvisc | 0.0000257 | Paxs | 905.26 | Joback Method |
| dvisc | 0.0000345 | Paxs | 837.63 | Joback Method |

| | | | | |
|-------|-----------|------|--------|---------------|
| dvisc | 0.0000486 | Paxs | 770.00 | Joback Method |
| dvisc | 0.0000733 | Paxs | 702.37 | Joback Method |
| dvisc | 0.0001206 | Paxs | 634.74 | Joback Method |
| dvisc | 0.0002235 | Paxs | 567.11 | Joback Method |
| dvisc | 0.0004895 | Paxs | 499.48 | Joback Method |

Sources

| | |
|------------------------|---|
| McGowan Method: | http://link.springer.com/article/10.1007/BF02311772 |
| NIST Webbook: | http://webbook.nist.gov/cgi/cbook.cgi?ID=U354312&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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