

Sebacic acid, di(2-heptyl) ester

Inchi: InChI=1S/C24H46O4/c1-5-7-13-17-21(3)27-23(25)19-15-11-9-10-12-16-20-24(26)28-22(29)
InchiKey: ZACMHOIXXJSCRZ-UHFFFAOYSA-N
Formula: C24H46O4
SMILES: CCCCCC(C)OC(=O)CCCCCCCCC(=O)OC(C)CCCC
Mol. weight [g/mol]: 398.62

Physical Properties

| Property code | Value | Unit | Source |
|---------------|----------|----------------------|----------------|
| gf | -321.52 | kJ/mol | Joback Method |
| hf | -1038.85 | kJ/mol | Joback Method |
| hfus | 56.44 | kJ/mol | Joback Method |
| hvap | 86.55 | kJ/mol | Joback Method |
| log10ws | -7.82 | | Crippen Method |
| logp | 7.131 | | Crippen Method |
| mvol | 363.900 | ml/mol | McGowan Method |
| pc | 857.97 | kPa | Joback Method |
| rinpol | 2603.00 | | NIST Webbook |
| tb | 900.22 | K | Joback Method |
| tc | 1102.59 | K | Joback Method |
| tf | 474.56 | K | Joback Method |
| vc | 1.415 | m ³ /kmol | Joback Method |

Temperature Dependent Properties

| Property code | Value | Unit | Temperature [K] | Source |
|---------------|-----------|---------|-----------------|---------------|
| cpg | 1218.42 | J/molxK | 900.22 | Joback Method |
| cpg | 1238.45 | J/molxK | 933.95 | Joback Method |
| cpg | 1257.07 | J/molxK | 967.68 | Joback Method |
| cpg | 1274.32 | J/molxK | 1001.41 | Joback Method |
| cpg | 1290.22 | J/molxK | 1035.14 | Joback Method |
| cpg | 1304.81 | J/molxK | 1068.86 | Joback Method |
| cpg | 1318.12 | J/molxK | 1102.59 | Joback Method |
| dvisc | 0.0007142 | Paxs | 474.56 | Joback Method |
| dvisc | 0.0002826 | Paxs | 545.50 | Joback Method |

| | | | | |
|-------|-----------|------|--------|---------------|
| dvisc | 0.0001385 | Paxs | 616.45 | Joback Method |
| dvisc | 0.0000786 | Paxs | 687.39 | Joback Method |
| dvisc | 0.0000496 | Paxs | 758.33 | Joback Method |
| dvisc | 0.0000339 | Paxs | 829.28 | Joback Method |
| dvisc | 0.0000246 | Paxs | 900.22 | Joback Method |

Sources

| | |
|------------------------|---|
| 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 |
| NIST Webbook: | http://webbook.nist.gov/cgi/cbook.cgi?ID=U355556&Units=SI |
| Crippen Method: | http://pubs.acs.org/doi/abs/10.1021/ci990307l |

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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<https://www.chemeo.com/cid/15-986-1/Sebacic-acid-di-2-heptyl-ester.pdf>

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