

Sebacic acid, dodecyl 3-oxobut-2-yl ester

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| Inchi: | InChI=1S/C26H48O5/c1-4-5-6-7-8-9-10-13-16-19-22-30-25(28)20-17-14-11-12-15-18-21 |
| InchiKey: | IOOTYRJIUMJVHY-UHFFFAOYSA-N |
| Formula: | C26H48O5 |
| SMILES: | CCCCCCCCCCCCOC(=O)CCCCCCCCC(=O)OC(C)C(C)=O |
| Mol. weight [g/mol]: | 440.66 |

Physical Properties

| Property code | Value | Unit | Source |
|---------------|----------|----------------------|----------------|
| gf | -431.16 | kJ/mol | Joback Method |
| hf | -1187.43 | kJ/mol | Joback Method |
| hfus | 66.75 | kJ/mol | Joback Method |
| hvap | 98.14 | kJ/mol | Joback Method |
| log10ws | -7.82 | | Crippen Method |
| logp | 7.092 | | Crippen Method |
| mcvol | 393.650 | ml/mol | McGowan Method |
| pc | 793.05 | kPa | Joback Method |
| rinpola | 3060.00 | | NIST Webbook |
| tb | 1000.29 | K | Joback Method |
| tc | 1235.03 | K | Joback Method |
| tf | 562.03 | K | Joback Method |
| vc | 1.540 | m ³ /kmol | Joback Method |

Temperature Dependent Properties

| Property code | Value | Unit | Temperature [K] | Source |
|---------------|-----------|---------|-----------------|---------------|
| cpg | 1365.44 | J/molxK | 1000.29 | Joback Method |
| cpg | 1444.26 | J/molxK | 1195.90 | Joback Method |
| cpg | 1432.02 | J/molxK | 1156.78 | Joback Method |
| cpg | 1418.08 | J/molxK | 1117.66 | Joback Method |
| cpg | 1402.37 | J/molxK | 1078.54 | Joback Method |
| cpg | 1384.85 | J/molxK | 1039.41 | Joback Method |
| cpg | 1454.85 | J/molxK | 1235.03 | Joback Method |
| dvisc | 0.0000183 | Paxs | 1000.29 | Joback Method |
| dvisc | 0.0000246 | Paxs | 927.25 | Joback Method |

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|-------|-----------|------|--------|---------------|
| dvisc | 0.0000349 | Paxs | 854.20 | Joback Method |
| dvisc | 0.0000528 | Paxs | 781.16 | Joback Method |
| dvisc | 0.0000870 | Paxs | 708.12 | Joback Method |
| dvisc | 0.0001607 | Paxs | 635.07 | Joback Method |
| dvisc | 0.0003483 | Paxs | 562.03 | Joback Method |

Sources

| | |
|------------------------|---|
| 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 |
| NIST Webbook: | http://webbook.nist.gov/cgi/cbook.cgi?ID=U355784&Units=SI |

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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