

Fumaric acid, 2-butyl octadecyl ester

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|-----------------------------|--|
| Inchi: | InChI=1S/C26H48O4/c1-4-6-7-8-9-10-11-12-13-14-15-16-17-18-19-20-23-29-25(27)21-2 |
| InchiKey: | YFACVHBJUIKMPE-QURGRASLSA-N |
| Formula: | C26H48O4 |
| SMILES: | CCCCCCCCCCCCCCCCCOC(=O)C=CC(=O)OC(C)CC |
| Mol. weight [g/mol]: | 424.66 |

Physical Properties

| Property code | Value | Unit | Source |
|---------------|---------|---------|----------------|
| gf | -222.02 | kJ/mol | Joback Method |
| hf | -957.63 | kJ/mol | Joback Method |
| hfus | 65.35 | kJ/mol | Joback Method |
| hvap | 91.35 | kJ/mol | Joback Method |
| log10ws | -8.40 | | Crippen Method |
| logp | 7.689 | | Crippen Method |
| mcvol | 387.780 | ml/mol | McGowan Method |
| pc | 788.16 | kPa | Joback Method |
| rinpol | 2931.00 | | NIST Webbook |
| tb | 950.58 | K | Joback Method |
| tc | 1167.61 | K | Joback Method |
| tf | 507.02 | K | Joback Method |
| vc | 1.514 | m3/kmol | Joback Method |

Temperature Dependent Properties

| Property code | Value | Unit | Temperature [K] | Source |
|---------------|-----------|---------|-----------------|---------------|
| cpg | 1317.93 | J/molxK | 950.58 | Joback Method |
| cpg | 1338.52 | J/molxK | 986.75 | Joback Method |
| cpg | 1357.62 | J/molxK | 1022.92 | Joback Method |
| cpg | 1375.27 | J/molxK | 1059.10 | Joback Method |
| cpg | 1391.54 | J/molxK | 1095.27 | Joback Method |
| cpg | 1406.49 | J/molxK | 1131.44 | Joback Method |
| cpg | 1420.19 | J/molxK | 1167.61 | Joback Method |
| dvisc | 0.0004362 | Paxs | 507.02 | Joback Method |
| dvisc | 0.0001808 | Paxs | 580.95 | Joback Method |

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|-------|-----------|------|--------|---------------|
| dvisc | 0.0000914 | Paxs | 654.87 | Joback Method |
| dvisc | 0.0000531 | Paxs | 728.80 | Joback Method |
| dvisc | 0.0000341 | Paxs | 802.73 | Joback Method |
| dvisc | 0.0000236 | Paxs | 876.65 | Joback Method |
| dvisc | 0.0000173 | Paxs | 950.58 | Joback Method |

Sources

| | |
|------------------------|---|
| McGowan Method: | http://link.springer.com/article/10.1007/BF02311772 |
| NIST Webbook: | http://webbook.nist.gov/cgi/cbook.cgi?ID=U348655&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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