

Fumaric acid, dec-4-enyl undecyl ester

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|-----------------------------|--|
| Inchi: | InChI=1S/C25H44O4/c1-3-5-7-9-11-13-15-17-19-23-29-25(27)21-20-24(26)28-22-18-16- |
| InchiKey: | ITWJPALOVIIXML-KFSAMSKSSA-N |
| Formula: | C25H44O4 |
| SMILES: | CCCCC=CCCCOC(=O)C=CC(=O)OCCCCCCCCCCC |
| Mol. weight [g/mol]: | 408.61 |

Physical Properties

| Property code | Value | Unit | Source |
|---------------|---------|----------------------|----------------|
| gf | -147.78 | kJ/mol | Joback Method |
| hf | -814.49 | kJ/mol | Joback Method |
| hfus | 66.48 | kJ/mol | Joback Method |
| hvap | 89.47 | kJ/mol | Joback Method |
| log10ws | -7.72 | | Crippen Method |
| logp | 7.076 | | Crippen Method |
| mvol | 369.390 | ml/mol | McGowan Method |
| pc | 853.96 | kPa | Joback Method |
| rinpol | 2864.00 | | NIST Webbook |
| rinpol | 2864.00 | | NIST Webbook |
| tb | 932.30 | K | Joback Method |
| tc | 1142.51 | K | Joback Method |
| tf | 505.67 | K | Joback Method |
| vc | 1.444 | m ³ /kmol | Joback Method |

Temperature Dependent Properties

| Property code | Value | Unit | Temperature [K] | Source |
|---------------|-----------|---------|-----------------|---------------|
| cpg | 1226.29 | J/molxK | 932.30 | Joback Method |
| cpg | 1312.29 | J/molxK | 1107.47 | Joback Method |
| cpg | 1297.32 | J/molxK | 1072.44 | Joback Method |
| cpg | 1281.31 | J/molxK | 1037.40 | Joback Method |
| cpg | 1264.18 | J/molxK | 1002.37 | Joback Method |
| cpg | 1245.86 | J/molxK | 967.33 | Joback Method |
| cpg | 1326.26 | J/molxK | 1142.51 | Joback Method |
| dvisc | 0.0000193 | Paxs | 932.30 | Joback Method |

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|-------|-----------|------|--------|---------------|
| dvisc | 0.0000260 | Paxs | 861.20 | Joback Method |
| dvisc | 0.0000370 | Paxs | 790.09 | Joback Method |
| dvisc | 0.0000562 | Paxs | 718.99 | Joback Method |
| dvisc | 0.0000938 | Paxs | 647.88 | Joback Method |
| dvisc | 0.0001776 | Paxs | 576.77 | Joback Method |
| dvisc | 0.0004022 | Paxs | 505.67 | Joback Method |

Sources

| | |
|------------------------|---|
| Crippen Method: | http://pubs.acs.org/doi/abs/10.1021/ci9903071 |
| Crippen Method: | https://www.cheméo.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=U348945&Units=SI |

Legend

| | |
|----------------------------|---|
| cp_g: | Ideal gas heat capacity |
| dvisc: | Dynamic viscosity |
| g_f: | Standard Gibbs free energy of formation |
| h_f: | Enthalpy of formation at standard conditions |
| h_{fus}: | Enthalpy of fusion at standard conditions |
| h_{vap}: | Enthalpy of vaporization at standard conditions |
| log₁₀ws: | Log ₁₀ of Water solubility in mol/l |
| log_p: | 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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