

Adipic acid, 3-ethylphenyl pentyl ester

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|-----------------------------|---|
| Inchi: | InChI=1S/C19H28O4/c1-3-5-8-14-22-18(20)12-6-7-13-19(21)23-17-11-9-10-16(4-2)15-17 |
| InchiKey: | TWMQSVPTDRFGHW-UHFFFAOYSA-N |
| Formula: | C19H28O4 |
| SMILES: | CCCCCOC(=O)CCCCC(=O)Oc1cccc(CC)c1 |
| Mol. weight [g/mol]: | 320.42 |

Physical Properties

| Property code | Value | Unit | Source |
|---------------|---------|----------------------|----------------|
| gf | -255.96 | kJ/mol | Joback Method |
| hf | -700.03 | kJ/mol | Joback Method |
| hfus | 44.19 | kJ/mol | Joback Method |
| hvap | 79.14 | kJ/mol | Joback Method |
| log10ws | -5.22 | | Crippen Method |
| logp | 4.448 | | Crippen Method |
| mvol | 269.690 | ml/mol | McGowan Method |
| pc | 1441.35 | kPa | Joback Method |
| rinpol | 2343.00 | | NIST Webbook |
| rinpol | 2343.00 | | NIST Webbook |
| tb | 818.36 | K | Joback Method |
| tc | 1017.87 | K | Joback Method |
| tf | 487.15 | K | Joback Method |
| vc | 1.040 | m ³ /kmol | Joback Method |

Temperature Dependent Properties

| Property code | Value | Unit | Temperature [K] | Source |
|---------------|-----------|---------|-----------------|---------------|
| cpg | 819.06 | J/molxK | 818.36 | Joback Method |
| cpg | 834.97 | J/molxK | 851.61 | Joback Method |
| cpg | 849.80 | J/molxK | 884.86 | Joback Method |
| cpg | 863.56 | J/molxK | 918.12 | Joback Method |
| cpg | 876.27 | J/molxK | 951.37 | Joback Method |
| cpg | 887.96 | J/molxK | 984.62 | Joback Method |
| cpg | 898.65 | J/molxK | 1017.87 | Joback Method |
| dvisc | 0.0006541 | Paxs | 487.15 | Joback Method |

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|-------|-----------|------|--------|---------------|
| dvisc | 0.0003630 | Paxs | 542.35 | Joback Method |
| dvisc | 0.0002246 | Paxs | 597.55 | Joback Method |
| dvisc | 0.0001507 | Paxs | 652.75 | Joback Method |
| dvisc | 0.0001077 | Paxs | 707.96 | Joback Method |
| dvisc | 0.0000807 | Paxs | 763.16 | Joback Method |
| dvisc | 0.0000629 | Paxs | 818.36 | 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=U353894&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/57-530-0/Adipic-acid-3-ethylphenyl-pentyl-ester.pdf>

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