

4-Ethylbenzoic acid, 4-hexadecyl ester

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|-----------------------------|--|
| Inchi: | InChI=1S/C25H42O2/c1-4-7-8-9-10-11-12-13-14-15-17-24(16-5-2)27-25(26)23-20-18-22 |
| InchiKey: | JLUAINKQSUGVNL-UHFFFAOYSA-N |
| Formula: | C25H42O2 |
| SMILES: | CCCCCCCCCCCC(CCC)OC(=O)c1ccc(CC)cc1 |
| Mol. weight [g/mol]: | 374.60 |

Physical Properties

| Property code | Value | Unit | Source |
|---------------|---------|----------------------|----------------|
| gf | 26.04 | kJ/mol | Joback Method |
| hf | -584.35 | kJ/mol | Joback Method |
| hfus | 53.42 | kJ/mol | Joback Method |
| hvap | 82.95 | kJ/mol | Joback Method |
| log10ws | -8.91 | | Crippen Method |
| logp | 7.886 | | Crippen Method |
| mvol | 346.790 | ml/mol | McGowan Method |
| pc | 947.91 | kPa | Joback Method |
| rinpol | 2604.70 | | NIST Webbook |
| rinpol | 2604.70 | | NIST Webbook |
| tb | 878.91 | K | Joback Method |
| tc | 1078.47 | K | Joback Method |
| tf | 467.61 | K | Joback Method |
| vc | 1.345 | m ³ /kmol | Joback Method |

Temperature Dependent Properties

| Property code | Value | Unit | Temperature [K] | Source |
|---------------|-----------|---------|-----------------|---------------|
| cpg | 1130.24 | J/molxK | 878.91 | Joback Method |
| cpg | 1149.81 | J/molxK | 912.17 | Joback Method |
| cpg | 1168.12 | J/molxK | 945.43 | Joback Method |
| cpg | 1185.24 | J/molxK | 978.69 | Joback Method |
| cpg | 1201.20 | J/molxK | 1011.95 | Joback Method |
| cpg | 1216.06 | J/molxK | 1045.21 | Joback Method |
| cpg | 1229.87 | J/molxK | 1078.47 | Joback Method |
| dvisc | 0.0007290 | Paxs | 467.61 | Joback Method |

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|-------|-----------|------|--------|---------------|
| dvisc | 0.0003164 | Paxs | 536.16 | Joback Method |
| dvisc | 0.0001659 | Paxs | 604.71 | Joback Method |
| dvisc | 0.0000992 | Paxs | 673.26 | Joback Method |
| dvisc | 0.0000653 | Paxs | 741.81 | Joback Method |
| dvisc | 0.0000461 | Paxs | 810.36 | Joback Method |
| dvisc | 0.0000344 | Paxs | 878.91 | 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=U292203&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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