

Isophthalic acid, 2-ethylphenyl heptyl ester

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|-----------------------------|--|
| Inchi: | InChI=1S/C23H28O4/c1-3-5-6-7-10-16-26-22(24)19-13-11-14-20(17-19)23(25)27-21-15- |
| InchiKey: | MKCKAFLMSWMCNQ-UHFFFAOYSA-N |
| Formula: | C23H28O4 |
| SMILES: | CCCCCCCOC(=O)c1cccc(C(=O)Oc2ccccc2CC)c1 |
| Mol. weight [g/mol]: | 368.47 |

Physical Properties

| Property code | Value | Unit | Source |
|---------------|---------|---------|----------------|
| gf | -119.50 | kJ/mol | Joback Method |
| hf | -557.53 | kJ/mol | Joback Method |
| hfus | 48.20 | kJ/mol | Joback Method |
| hvap | 90.98 | kJ/mol | Joback Method |
| log10ws | -7.12 | | Crippen Method |
| logp | 5.595 | | Crippen Method |
| mcvol | 302.290 | ml/mol | McGowan Method |
| pc | 1369.71 | kPa | Joback Method |
| rinpol | 2805.00 | | NIST Webbook |
| tb | 941.54 | K | Joback Method |
| tc | 1164.15 | K | Joback Method |
| tf | 571.17 | K | Joback Method |
| vc | 1.155 | m3/kmol | Joback Method |

Temperature Dependent Properties

| Property code | Value | Unit | Temperature [K] | Source |
|---------------|-----------|---------|-----------------|---------------|
| cpg | 955.96 | J/molxK | 941.54 | Joback Method |
| cpg | 970.23 | J/molxK | 978.64 | Joback Method |
| cpg | 983.14 | J/molxK | 1015.74 | Joback Method |
| cpg | 994.72 | J/molxK | 1052.85 | Joback Method |
| cpg | 1005.01 | J/molxK | 1089.95 | Joback Method |
| cpg | 1014.06 | J/molxK | 1127.05 | Joback Method |
| cpg | 1021.91 | J/molxK | 1164.15 | Joback Method |
| dvisc | 0.0003350 | Paxs | 571.17 | Joback Method |
| dvisc | 0.0001947 | Paxs | 632.90 | Joback Method |

| | | | | |
|-------|-----------|------|--------|---------------|
| dvisc | 0.0001247 | Paxs | 694.63 | Joback Method |
| dvisc | 0.0000858 | Paxs | 756.36 | Joback Method |
| dvisc | 0.0000625 | Paxs | 818.08 | Joback Method |
| dvisc | 0.0000476 | Paxs | 879.81 | Joback Method |
| dvisc | 0.0000376 | Paxs | 941.54 | Joback Method |

Sources

| | |
|------------------------|---|
| McGowan Method: | http://link.springer.com/article/10.1007/BF02311772 |
| NIST Webbook: | http://webbook.nist.gov/cgi/cbook.cgi?ID=U356765&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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