

Isophthalic acid, dodecyl 4-methylpent-2-yl ester

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|----------------------|--|
| Inchi: | InChI=1S/C26H42O4/c1-5-6-7-8-9-10-11-12-13-14-18-29-25(27)23-16-15-17-24(20-23)2 |
| InchiKey: | XPHARKGMSUBWIU-UHFFFAOYSA-N |
| Formula: | C26H42O4 |
| SMILES: | CCCCCCCCCCCCOC(=O)c1cccc(C(=O)OC(C)CC(C)C)c1 |
| Mol. weight [g/mol]: | 418.61 |

Physical Properties

| Property code | Value | Unit | Source |
|---------------|---------|----------------------|----------------|
| gf | -201.90 | kJ/mol | Joback Method |
| hf | -855.07 | kJ/mol | Joback Method |
| hfus | 55.28 | kJ/mol | Joback Method |
| hvap | 93.94 | kJ/mol | Joback Method |
| log10ws | -8.52 | | Crippen Method |
| logp | 7.356 | | Crippen Method |
| mvol | 368.320 | ml/mol | McGowan Method |
| pc | 923.86 | kPa | Joback Method |
| rinpol | 2899.00 | | NIST Webbook |
| tb | 977.64 | K | Joback Method |
| tc | 1196.94 | K | Joback Method |
| tf | 536.04 | K | Joback Method |
| vc | 1.419 | m ³ /kmol | Joback Method |

Temperature Dependent Properties

| Property code | Value | Unit | Temperature [K] | Source |
|---------------|-----------|---------|-----------------|---------------|
| cpg | 1244.40 | J/molxK | 977.64 | Joback Method |
| cpg | 1316.39 | J/molxK | 1160.39 | Joback Method |
| cpg | 1304.91 | J/molxK | 1123.84 | Joback Method |
| cpg | 1292.02 | J/molxK | 1087.29 | Joback Method |
| cpg | 1277.68 | J/molxK | 1050.74 | Joback Method |
| cpg | 1261.82 | J/molxK | 1014.19 | Joback Method |
| cpg | 1326.50 | J/molxK | 1196.94 | Joback Method |
| dvisc | 0.0000189 | Paxs | 977.64 | Joback Method |
| dvisc | 0.0000254 | Paxs | 904.04 | Joback Method |

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|-------|-----------|------|--------|---------------|
| dvisc | 0.0000360 | Paxs | 830.44 | Joback Method |
| dvisc | 0.0000545 | Paxs | 756.84 | Joback Method |
| dvisc | 0.0000904 | Paxs | 683.24 | Joback Method |
| dvisc | 0.0001694 | Paxs | 609.64 | Joback Method |
| dvisc | 0.0003771 | Paxs | 536.04 | 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=U356449&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/19-639-2/Isophthalic-acid-dodecyl-4-methylpent-2-yl-ester.pdf>

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