

Succinic acid, dodecyl 2,2,2-trichloroethyl ester

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|----------------------|--|
| Inchi: | InChI=1S/C18H31Cl3O4/c1-2-3-4-5-6-7-8-9-10-11-14-24-16(22)12-13-17(23)25-15-18(19) |
| InchiKey: | BBLDVAZHTUMKSA-UHFFFAOYSA-N |
| Formula: | C18H31Cl3O4 |
| SMILES: | CCCCCCCCCCCCOC(=O)CCC(=O)OCC(Cl)(Cl)Cl |
| Mol. weight [g/mol]: | 417.80 |

Physical Properties

| Property code | Value | Unit | Source |
|---------------|---------|----------------------|----------------|
| gf | -400.11 | kJ/mol | Joback Method |
| hf | -960.42 | kJ/mol | Joback Method |
| hfus | 53.13 | kJ/mol | Joback Method |
| hvap | 85.83 | kJ/mol | Joback Method |
| log10ws | -6.64 | | Crippen Method |
| logp | 6.144 | | Crippen Method |
| mvol | 316.080 | ml/mol | McGowan Method |
| pc | 1152.22 | kPa | Joback Method |
| rinpol | 2568.00 | | NIST Webbook |
| tb | 872.88 | K | Joback Method |
| tc | 1072.77 | K | Joback Method |
| tf | 529.12 | K | Joback Method |
| vc | 1.228 | m ³ /kmol | Joback Method |

Temperature Dependent Properties

| Property code | Value | Unit | Temperature [K] | Source |
|---------------|-----------|---------|-----------------|---------------|
| cpg | 938.85 | J/molxK | 872.88 | Joback Method |
| cpg | 953.46 | J/molxK | 906.19 | Joback Method |
| cpg | 967.05 | J/molxK | 939.51 | Joback Method |
| cpg | 979.68 | J/molxK | 972.82 | Joback Method |
| cpg | 991.37 | J/molxK | 1006.14 | Joback Method |
| cpg | 1002.17 | J/molxK | 1039.45 | Joback Method |
| cpg | 1012.11 | J/molxK | 1072.77 | Joback Method |
| dvisc | 0.0004661 | Paxs | 529.12 | Joback Method |
| dvisc | 0.0002444 | Paxs | 586.41 | Joback Method |

| | | | | |
|-------|-----------|------|--------|---------------|
| dvisc | 0.0001438 | Paxs | 643.71 | Joback Method |
| dvisc | 0.0000922 | Paxs | 701.00 | Joback Method |
| dvisc | 0.0000633 | Paxs | 758.29 | Joback Method |
| dvisc | 0.0000458 | Paxs | 815.59 | Joback Method |
| dvisc | 0.0000345 | Paxs | 872.88 | Joback Method |

Sources

| | |
|------------------------|---|
| McGowan Method: | http://link.springer.com/article/10.1007/BF02311772 |
| NIST Webbook: | http://webbook.nist.gov/cgi/cbook.cgi?ID=U349175&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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