

3-Octenoic acid, undecyl ester

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|-----------------------------|--|
| Inchi: | InChI=1S/C19H36O2/c1-3-5-7-9-10-11-12-14-16-18-21-19(20)17-15-13-8-6-4-2/h13,15H |
| InchiKey: | MNPJNTQGBSAOGZ-FYWRMAATSA-N |
| Formula: | C19H36O2 |
| SMILES: | CCCCC=CCC(=O)OCCCCCCCCCCC |
| Mol. weight [g/mol]: | 296.49 |

Physical Properties

| Property code | Value | Unit | Source |
|---------------|---------|----------------------|----------------|
| gf | -44.60 | kJ/mol | Joback Method |
| hf | -563.07 | kJ/mol | Joback Method |
| hfus | 47.95 | kJ/mol | Joback Method |
| hvap | 67.00 | kJ/mol | Joback Method |
| log10ws | -6.49 | | Crippen Method |
| logp | 6.197 | | Crippen Method |
| mvol | 281.710 | ml/mol | McGowan Method |
| pc | 1153.78 | kPa | Joback Method |
| rinpol | 2080.00 | | NIST Webbook |
| rinpol | 2080.00 | | NIST Webbook |
| tb | 714.57 | K | Joback Method |
| tc | 887.77 | K | Joback Method |
| tf | 370.97 | K | Joback Method |
| vc | 1.103 | m ³ /kmol | Joback Method |

Temperature Dependent Properties

| Property code | Value | Unit | Temperature [K] | Source |
|---------------|-----------|---------|-----------------|---------------|
| cpg | 826.19 | J/molxK | 714.57 | Joback Method |
| cpg | 911.43 | J/molxK | 858.90 | Joback Method |
| cpg | 896.00 | J/molxK | 830.04 | Joback Method |
| cpg | 879.80 | J/molxK | 801.17 | Joback Method |
| cpg | 862.78 | J/molxK | 772.30 | Joback Method |
| cpg | 844.92 | J/molxK | 743.44 | Joback Method |
| cpg | 926.09 | J/molxK | 887.77 | Joback Method |
| dvisc | 0.0000724 | Paxs | 714.57 | Joback Method |

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|-------|-----------|------|--------|---------------|
| dvisc | 0.0000975 | Paxs | 657.30 | Joback Method |
| dvisc | 0.0001390 | Paxs | 600.04 | Joback Method |
| dvisc | 0.0002137 | Paxs | 542.77 | Joback Method |
| dvisc | 0.0003634 | Paxs | 485.50 | Joback Method |
| dvisc | 0.0007124 | Paxs | 428.24 | Joback Method |
| dvisc | 0.0017193 | Paxs | 370.97 | Joback Method |

Sources

| | |
|------------------------|---|
| 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=U406131&Units=SI |
| Crippen Method: | http://pubs.acs.org/doi/abs/10.1021/ci990307l |
| Crippen Method: | https://www.chemeo.com/doc/models/crippen_log10ws |

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/89-421-6/3-Octenoic-acid-undecyl-ester.pdf>

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