

Oct-3-enoic acid, undec-2-en-1-yl ester

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|-----------------------------|--|
| Inchi: | InChI=1S/C19H34O2/c1-3-5-7-9-10-11-12-14-16-18-21-19(20)17-15-13-8-6-4-2/h13-16H |
| InchiKey: | PBZKKWCQDPDQII-WXUKJITCSA-N |
| Formula: | C19H34O2 |
| SMILES: | CCCCC=CCC(=O)OCC=CCCCCCCCC |
| Mol. weight [g/mol]: | 294.47 |

Physical Properties

| Property code | Value | Unit | Source |
|---------------|---------|----------------------|----------------|
| gf | 35.62 | kJ/mol | Joback Method |
| hf | -445.85 | kJ/mol | Joback Method |
| hfus | 48.16 | kJ/mol | Joback Method |
| hvap | 66.96 | kJ/mol | Joback Method |
| log10ws | -6.35 | | Crippen Method |
| logp | 5.973 | | Crippen Method |
| mcvol | 277.410 | ml/mol | McGowan Method |
| pc | 1195.65 | kPa | Joback Method |
| rinpol | 2082.00 | | NIST Webbook |
| rinpol | 2082.00 | | NIST Webbook |
| tb | 718.73 | K | Joback Method |
| tc | 896.01 | K | Joback Method |
| tf | 365.89 | K | Joback Method |
| vc | 1.083 | m ³ /kmol | Joback Method |

Temperature Dependent Properties

| Property code | Value | Unit | Temperature [K] | Source |
|---------------|-----------|---------|-----------------|---------------|
| cpg | 803.45 | J/molxK | 718.73 | Joback Method |
| cpg | 886.41 | J/molxK | 866.47 | Joback Method |
| cpg | 871.40 | J/molxK | 836.92 | Joback Method |
| cpg | 855.63 | J/molxK | 807.37 | Joback Method |
| cpg | 839.08 | J/molxK | 777.82 | Joback Method |
| cpg | 821.70 | J/molxK | 748.28 | Joback Method |
| cpg | 900.71 | J/molxK | 896.01 | Joback Method |
| dvisc | 0.0000626 | Paxs | 718.73 | Joback Method |

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|-------|-----------|------|--------|---------------|
| dvisc | 0.0000845 | Paxs | 659.92 | Joback Method |
| dvisc | 0.0001209 | Paxs | 601.12 | Joback Method |
| dvisc | 0.0001871 | Paxs | 542.31 | Joback Method |
| dvisc | 0.0003219 | Paxs | 483.50 | Joback Method |
| dvisc | 0.0006439 | Paxs | 424.70 | Joback Method |
| dvisc | 0.0016090 | Paxs | 365.89 | Joback Method |

Sources

| | |
|------------------------|---|
| Crippen Method: | http://pubs.acs.org/doi/abs/10.1021/ci9903071 |
| Crippen Method: | https://www.cheméo.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=U406953&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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