

Octadeca-9,12-dienoic acid octadeca-9,12-dienyl ester, Z,Z,Z,Z

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|-----------------------------|--|
| Inchi: | InChI=1S/C36H64O2/c1-3-5-7-9-11-13-15-17-19-21-23-25-27-29-31-33-35-38-36(37)34- |
| InchiKey: | BWPODKUVJMQURN-YIEDHNAISA-N |
| Formula: | C36H64O2 |
| SMILES: | CCCCC=CCC=CCCCCCCCCOC(=O)CCCCCCCC=CCC=CCCCC |
| Mol. weight [g/mol]: | 528.89 |

Physical Properties

| Property code | Value | Unit | Source |
|---------------|---------|----------------------|----------------|
| gf | 339.20 | kJ/mol | Joback Method |
| hf | -562.29 | kJ/mol | Joback Method |
| hfus | 92.59 | kJ/mol | Joback Method |
| hvap | 104.72 | kJ/mol | Joback Method |
| log10ws | -13.17 | | Crippen Method |
| logp | 12.157 | | Crippen Method |
| mvol | 508.340 | ml/mol | McGowan Method |
| pc | 511.87 | kPa | Joback Method |
| rinpol | 3685.73 | | NIST Webbook |
| rinpol | 3685.73 | | NIST Webbook |
| tb | 1116.01 | K | Joback Method |
| tc | 1420.29 | K | Joback Method |
| tf | 547.32 | K | Joback Method |
| vc | 1.996 | m ³ /kmol | Joback Method |

Temperature Dependent Properties

| Property code | Value | Unit | Temperature [K] | Source |
|---------------|-----------|---------|-----------------|---------------|
| cpg | 1844.08 | J/molxK | 1116.01 | Joback Method |
| cpg | 2000.75 | J/molxK | 1369.58 | Joback Method |
| cpg | 1970.18 | J/molxK | 1318.87 | Joback Method |
| cpg | 1939.65 | J/molxK | 1268.15 | Joback Method |
| cpg | 1908.75 | J/molxK | 1217.44 | Joback Method |
| cpg | 1877.03 | J/molxK | 1166.72 | Joback Method |
| cpg | 2031.78 | J/molxK | 1420.29 | Joback Method |
| dvisc | 0.0000039 | Paxs | 1116.01 | Joback Method |

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|-------|-----------|------|---------|---------------|
| dvisc | 0.0000054 | Paxs | 1021.23 | Joback Method |
| dvisc | 0.0000081 | Paxs | 926.45 | Joback Method |
| dvisc | 0.0000132 | Paxs | 831.66 | Joback Method |
| dvisc | 0.0000243 | Paxs | 736.88 | Joback Method |
| dvisc | 0.0000539 | Paxs | 642.10 | Joback Method |
| dvisc | 0.0001575 | Paxs | 547.32 | Joback Method |

Sources

| | |
|------------------------|---|
| NIST Webbook: | http://webbook.nist.gov/cgi/cbook.cgi?ID=R437111&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 |
| McGowan Method: | http://link.springer.com/article/10.1007/BF02311772 |

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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