

# Cyclotridecene, 1-methoxy-, (Z)-

|                             |  |
|-----------------------------|--|
| <b>Inchi:</b>               | InChI=1S/C14H26O/c1-15-14-12-10-8-6-4-2-3-5-7-9-11-13-14/h12H,2-11,13H2,1H3/b14- |
| <b>InchiKey:</b>            | DQNSWVPJRZTUNF-WYMLVPIESA-N  |
| <b>Formula:</b>             | C14H26O  |
| <b>SMILES:</b>              | COC1=CCCCCCCCCCC1  |
| <b>Mol. weight [g/mol]:</b> | 210.36   |
| <b>CAS:</b>                 | 78289-14-2   |

## Physical Properties

| Property code | Value   | Unit    | Source         |
|---------------|---------|---------|----------------|
| gf            | -70.21  | kJ/mol  | Joback Method  |
| hf            | -386.66 | kJ/mol  | Joback Method  |
| hfus          | 10.10   | kJ/mol  | Joback Method  |
| hvap          | 52.06   | kJ/mol  | Joback Method  |
| log10ws       | -5.01   |         | Crippen Method |
| logp          | 4.821   |         | Crippen Method |
| mcvol         | 198.830 | ml/mol  | McGowan Method |
| pc            | 2173.43 | kPa     | Joback Method  |
| tb            | 600.39  | K       | Joback Method  |
| tc            | 839.73  | K       | Joback Method  |
| tf            | 270.03  | K       | Joback Method  |
| vc            | 0.702   | m3/kmol | Joback Method  |

## Temperature Dependent Properties

| Property code | Value     | Unit    | Temperature [K] | Source        |
|---------------|-----------|---------|-----------------|---------------|
| cpg           | 523.64    | J/molxK | 600.39          | Joback Method |
| cpg           | 551.00    | J/molxK | 640.28          | Joback Method |
| cpg           | 576.65    | J/molxK | 680.17          | Joback Method |
| cpg           | 600.57    | J/molxK | 720.06          | Joback Method |
| cpg           | 622.72    | J/molxK | 759.95          | Joback Method |
| cpg           | 643.06    | J/molxK | 799.84          | Joback Method |
| cpg           | 661.56    | J/molxK | 839.73          | Joback Method |
| dvisc         | 0.0252648 | Paxs    | 270.03          | Joback Method |
| dvisc         | 0.0028674 | Paxs    | 325.09          | Joback Method |

|       |           |      |        |               |
|-------|-----------|------|--------|---------------|
| dvisc | 0.0006112 | Paxs | 380.15 | Joback Method |
| dvisc | 0.0001927 | Paxs | 435.21 | Joback Method |
| dvisc | 0.0000787 | Paxs | 490.27 | Joback Method |
| dvisc | 0.0000385 | Paxs | 545.33 | Joback Method |
| dvisc | 0.0000215 | Paxs | 600.39 | Joback Method |

## Sources

|                        |   |
|------------------------|---|
| <b>Crippen Method:</b> | <a href="https://www.chemeo.com/doc/models/crippen_log10ws">https://www.chemeo.com/doc/models/crippen_log10ws</a>                             |
| <b>Joback Method:</b>  | <a href="https://en.wikipedia.org/wiki/Joback_method">https://en.wikipedia.org/wiki/Joback_method</a>   |
| <b>McGowan Method:</b> | <a href="http://link.springer.com/article/10.1007/BF02311772">http://link.springer.com/article/10.1007/BF02311772</a>                         |
| <b>NIST Webbook:</b>   | <a href="http://webbook.nist.gov/cgi/cbook.cgi?ID=C78289142&amp;Units=SI">http://webbook.nist.gov/cgi/cbook.cgi?ID=C78289142&amp;Units=SI</a> |
| <b>Crippen Method:</b> | <a href="http://pubs.acs.org/doi/abs/10.1021/ci990307l">http://pubs.acs.org/doi/abs/10.1021/ci990307l</a>                                     |

## Legend

|                 |   |
|-----------------|---|
| <b>cpg:</b>     | Ideal gas heat capacity                         |
| <b>dvisc:</b>   | Dynamic viscosity                               |
| <b>gf:</b>      | Standard Gibbs free energy of formation         |
| <b>hf:</b>      | Enthalpy of formation at standard conditions    |
| <b>hfus:</b>    | Enthalpy of fusion at standard conditions       |
| <b>hvap:</b>    | Enthalpy of vaporization at standard conditions |
| <b>log10ws:</b> | Log10 of Water solubility in mol/l              |
| <b>logp:</b>    | Octanol/Water partition coefficient             |
| <b>mcvol:</b>   | McGowan's characteristic volume                 |
| <b>pc:</b>      | Critical Pressure                               |
| <b>tb:</b>      | Normal Boiling Point Temperature                |
| <b>tc:</b>      | Critical Temperature                            |
| <b>tf:</b>      | Normal melting (fusion) point                   |
| <b>vc:</b>      | Critical Volume                                 |

Latest version available from:

<https://www.chemeo.com/cid/16-133-6/Cyclotridecene-1-methoxy-Z.pdf>

Generated by Cheméo on 2024-04-27 05:00:57.42997461 +0000 UTC m=+16483306.350551925.

Cheméo (<https://www.chemeo.com>) is the biggest free database of chemical and physical data for the process industry.