

# (E,E,Z)-2,4,8-undecatrienal

|                      |  |
|----------------------|--|
| Inchi:               | InChI=1S/C11H16O/c1-2-3-4-5-6-7-8-9-10-11-12/h3-4,7-11H,2,5-6H2,1H3/b4-3-,8-7+,10- |
| InchiKey:            | JVVHINVPEGIIED-IICWQLAKSA-N  |
| Formula:             | C11H16O  |
| SMILES:              | CCC=CCCC=CC=CC=O   |
| Mol. weight [g/mol]: | 164.24   |

## Physical Properties

| Property code | Value   | Unit    | Source         |
|---------------|---------|---------|----------------|
| gf            | 182.88  | kJ/mol  | Joback Method  |
| hf            | -4.29   | kJ/mol  | Joback Method  |
| hfus          | 27.14   | kJ/mol  | Joback Method  |
| hvap          | 46.67   | kJ/mol  | Joback Method  |
| log10ws       | -3.27   |         | Crippen Method |
| logp          | 3.044   |         | Crippen Method |
| mcvol         | 154.520 | ml/mol  | McGowan Method |
| pc            | 2393.53 | kPa     | Joback Method  |
| ripol         | 1413.00 |         | NIST Webbook   |
| ripol         | 1965.00 |         | NIST Webbook   |
| ripol         | 1965.00 |         | NIST Webbook   |
| tb            | 512.22  | K       | Joback Method  |
| tc            | 702.68  | K       | Joback Method  |
| tf            | 240.49  | K       | Joback Method  |
| vc            | 0.609   | m3/kmol | Joback Method  |

## Temperature Dependent Properties

| Property code | Value  | Unit    | Temperature [K] | Source        |
|---------------|--------|---------|-----------------|---------------|
| cpg           | 336.68 | J/molxK | 512.22          | Joback Method |
| cpg           | 398.37 | J/molxK | 670.94          | Joback Method |
| cpg           | 387.42 | J/molxK | 639.19          | Joback Method |
| cpg           | 375.83 | J/molxK | 607.45          | Joback Method |
| cpg           | 363.54 | J/molxK | 575.71          | Joback Method |
| cpg           | 350.51 | J/molxK | 543.96          | Joback Method |
| cpg           | 408.72 | J/molxK | 702.68          | Joback Method |

|       |           |      |        |               |
|-------|-----------|------|--------|---------------|
| dvisc | 0.0001639 | Paxs | 512.22 | Joback Method |
| dvisc | 0.0002163 | Paxs | 466.93 | Joback Method |
| dvisc | 0.0003032 | Paxs | 421.64 | Joback Method |
| dvisc | 0.0004609 | Paxs | 376.36 | Joback Method |
| dvisc | 0.0007858 | Paxs | 331.07 | Joback Method |
| dvisc | 0.0015863 | Paxs | 285.78 | Joback Method |
| dvisc | 0.0041723 | Paxs | 240.49 | 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=R237020&amp;Units=SI">http://webbook.nist.gov/cgi/cbook.cgi?ID=R237020&amp;Units=SI</a> |
| <b>Crippen Method:</b> | <a href="http://pubs.acs.org/doi/abs/10.1021/ci9903071">http://pubs.acs.org/doi/abs/10.1021/ci9903071</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>rinpol:</b>  | Non-polar retention indices                     |
| <b>ripol:</b>   | Polar retention indices                         |
| <b>tb:</b>      | Normal Boiling Point Temperature                |
| <b>tc:</b>      | Critical Temperature                            |
| <b>tf:</b>      | Normal melting (fusion) point                   |
| <b>vc:</b>      | Critical Volume                                 |

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