

# 7-dodecenal, E

|                      |                                                                               |
|----------------------|-------------------------------------------------------------------------------|
| Other names:         | (E)-7-Dodecenal                                                               |
| Inchi:               | InChI=1S/C12H22O/c1-2-3-4-5-6-7-8-9-10-11-12-13/h5-6,12H,2-4,7-11H2,1H3/b6-5+ |
| InchiKey:            | HTUHYXDEKWCDCI-AATRIKPKSA-N                                                   |
| Formula:             | C12H22O                                                                       |
| SMILES:              | CCCCC=CCCCCCC=O                                                               |
| Mol. weight [g/mol]: | 182.30                                                                        |

## Physical Properties

| Property code | Value   | Unit    | Source         |
|---------------|---------|---------|----------------|
| gf            | 30.86   | kJ/mol  | Joback Method  |
| hf            | -259.37 | kJ/mol  | Joback Method  |
| hfus          | 29.33   | kJ/mol  | Joback Method  |
| hvap          | 48.98   | kJ/mol  | Joback Method  |
| log10ws       | -3.98   |         | Crippen Method |
| logp          | 3.882   |         | Crippen Method |
| mcvol         | 177.210 | ml/mol  | McGowan Method |
| pc            | 1985.89 | kPa     | Joback Method  |
| rinpol        | 1397.00 |         | NIST Webbook   |
| rinpol        | 1397.00 |         | NIST Webbook   |
| ripol         | 1732.00 |         | NIST Webbook   |
| tb            | 526.78  | K       | Joback Method  |
| tc            | 700.00  | K       | Joback Method  |
| tf            | 261.92  | K       | Joback Method  |
| vc            | 0.705   | m3/kmol | Joback Method  |

## Temperature Dependent Properties

| Property code | Value  | Unit    | Temperature [K] | Source        |
|---------------|--------|---------|-----------------|---------------|
| cpg           | 418.84 | J/mol×K | 526.78          | Joback Method |
| cpg           | 434.14 | J/mol×K | 555.65          | Joback Method |
| cpg           | 448.75 | J/mol×K | 584.52          | Joback Method |
| cpg           | 462.70 | J/mol×K | 613.39          | Joback Method |
| cpg           | 476.02 | J/mol×K | 642.26          | Joback Method |
| cpg           | 488.73 | J/mol×K | 671.13          | Joback Method |

|       |           |         |        |               |
|-------|-----------|---------|--------|---------------|
| cpg   | 500.86    | J/mol×K | 700.00 | Joback Method |
| dvisc | 0.0046993 | Pa×s    | 261.92 | Joback Method |
| dvisc | 0.0019373 | Pa×s    | 306.06 | Joback Method |
| dvisc | 0.0009985 | Pa×s    | 350.21 | Joback Method |
| dvisc | 0.0005970 | Pa×s    | 394.35 | Joback Method |
| dvisc | 0.0003959 | Pa×s    | 438.49 | Joback Method |
| dvisc | 0.0002830 | Pa×s    | 482.64 | Joback Method |
| dvisc | 0.0002140 | Pa×s    | 526.78 | Joback Method |

## Sources

|                        |                                                                                                                                           |
|------------------------|-------------------------------------------------------------------------------------------------------------------------------------------|
| <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=R265406&amp;Units=SI">http://webbook.nist.gov/cgi/cbook.cgi?ID=R265406&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>                                 |
| <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>                                     |

## 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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