

# 6-Octen-2-one, (Z)-

|                      |                                                            |
|----------------------|------------------------------------------------------------|
| Inchi:               | InChI=1S/C8H14O/c1-3-4-5-6-7-8(2)9/h3-4H,5-7H2,1-2H3/b4-3- |
| InchiKey:            | FPXPWFCLCGOFSTQ-ARJAWSKDSA-N                               |
| Formula:             | C8H14O                                                     |
| SMILES:              | CC=CCCCC(C)=O                                              |
| Mol. weight [g/mol]: | 126.20                                                     |
| CAS:                 | 74810-53-0                                                 |

## Physical Properties

| Property code | Value   | Unit    | Source         |
|---------------|---------|---------|----------------|
| gf            | -32.22  | kJ/mol  | Joback Method  |
| hf            | -203.81 | kJ/mol  | Joback Method  |
| hfus          | 18.28   | kJ/mol  | Joback Method  |
| hvap          | 40.11   | kJ/mol  | Joback Method  |
| log10ws       | -2.30   |         | Crippen Method |
| logp          | 2.322   |         | Crippen Method |
| mcvol         | 120.850 | ml/mol  | McGowan Method |
| pc            | 2856.62 | kPa     | Joback Method  |
| ripol         | 1316.00 |         | NIST Webbook   |
| tb            | 440.47  | K       | Joback Method  |
| tc            | 624.44  | K       | Joback Method  |
| tf            | 224.77  | K       | Joback Method  |
| vc            | 0.469   | m3/kmol | Joback Method  |

## Temperature Dependent Properties

| Property code | Value     | Unit    | Temperature [K] | Source        |
|---------------|-----------|---------|-----------------|---------------|
| cpg           | 239.73    | J/mol×K | 440.47          | Joback Method |
| cpg           | 295.21    | J/mol×K | 593.78          | Joback Method |
| cpg           | 285.17    | J/mol×K | 563.12          | Joback Method |
| cpg           | 274.62    | J/mol×K | 532.46          | Joback Method |
| cpg           | 263.55    | J/mol×K | 501.79          | Joback Method |
| cpg           | 251.92    | J/mol×K | 471.13          | Joback Method |
| cpg           | 304.76    | J/mol×K | 624.44          | Joback Method |
| dvisc         | 0.0002453 | Pa×s    | 440.47          | Joback Method |

|       |           |      |        |               |
|-------|-----------|------|--------|---------------|
| dvisc | 0.0003179 | Pa×s | 404.52 | Joback Method |
| dvisc | 0.0004335 | Pa×s | 368.57 | Joback Method |
| dvisc | 0.0006321 | Pa×s | 332.62 | Joback Method |
| dvisc | 0.0010100 | Pa×s | 296.67 | Joback Method |
| dvisc | 0.0018362 | Pa×s | 260.72 | Joback Method |
| dvisc | 0.0040418 | Pa×s | 224.77 | Joback Method |

## Sources

|                        |                                                                                                                                               |
|------------------------|-----------------------------------------------------------------------------------------------------------------------------------------------|
| <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=C74810530&amp;Units=SI">http://webbook.nist.gov/cgi/cbook.cgi?ID=C74810530&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>                             |

## 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>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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