

# Spiro[4.5]decan-2-one

|                             |   |
|-----------------------------|---|
| <b>Other names:</b>         | Spiro[4,5]decan-2-one                                 |
| <b>Inchi:</b>               | InChI=1S/C10H16O/c11-9-4-7-10(8-9)5-2-1-3-6-10/h1-8H2 |
| <b>InchiKey:</b>            | DJHRXBYWKVQDJY-UHFFFAOYSA-N                           |
| <b>Formula:</b>             | C10H16O   |
| <b>SMILES:</b>              | O=C1CCC2(CCCCC2)C1                                    |
| <b>Mol. weight [g/mol]:</b> | 152.23  |
| <b>CAS:</b>                 | 3643-16-1   |

## Physical Properties

| Property code | Value   | Unit                 | Source         |
|---------------|---------|----------------------|----------------|
| gf            | -13.95  | kJ/mol               | Joback Method  |
| hf            | -230.89 | kJ/mol               | Joback Method  |
| hfus          | 1.67    | kJ/mol               | Joback Method  |
| hvap          | 41.77   | kJ/mol               | Joback Method  |
| log10ws       | -2.84   |                      | Crippen Method |
| logp          | 2.690   |                      | Crippen Method |
| mcvol         | 131.610 | ml/mol               | McGowan Method |
| pc            | 3439.94 | kPa                  | Joback Method  |
| rinpol        | 1180.00 |                      | NIST Webbook   |
| tb            | 531.49  | K                    | Joback Method  |
| tc            | 781.63  | K                    | Joback Method  |
| tf            | 320.62  | K                    | Joback Method  |
| vc            | 0.483   | m <sup>3</sup> /kmol | Joback Method  |

## Temperature Dependent Properties

| Property code | Value  | Unit    | Temperature [K] | Source        |
|---------------|--------|---------|-----------------|---------------|
| cpg           | 323.99 | J/mol×K | 531.49          | Joback Method |
| cpg           | 344.73 | J/mol×K | 573.18          | Joback Method |
| cpg           | 363.97 | J/mol×K | 614.87          | Joback Method |
| cpg           | 381.88 | J/mol×K | 656.56          | Joback Method |
| cpg           | 398.65 | J/mol×K | 698.25          | Joback Method |
| cpg           | 414.43 | J/mol×K | 739.94          | Joback Method |
| cpg           | 429.42 | J/mol×K | 781.63          | Joback Method |

# Sources

|                        |   |
|------------------------|---|
| <b>NIST Webbook:</b>   | <a href="http://webbook.nist.gov/cgi/cbook.cgi?ID=C3643161&amp;Units=SI">http://webbook.nist.gov/cgi/cbook.cgi?ID=C3643161&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>                                   |
| <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>                       |

# Legend

|                 |   |
|-----------------|---|
| <b>cpg:</b>     | Ideal gas heat capacity                         |
| <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>hvp:</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>rinp:</b>    | Non-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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