

# p-Mentha-1,4-dien-9-al

|                             |   |
|-----------------------------|---|
| <b>Inchi:</b>               | InChI=1S/C10H14O/c1-8-3-5-10(6-4-8)9(2)7-11/h3,6-7,9H,4-5H2,1-2H3 |
| <b>InchiKey:</b>            | MWUMCBLYLOUAPN-UHFFFAOYSA-N                                       |
| <b>Formula:</b>             | C10H14O   |
| <b>SMILES:</b>              | CC1=CCC(C(C)C=O)=CC1  |
| <b>Mol. weight [g/mol]:</b> | 150.22  |

## Physical Properties

| Property code | Value   | Unit    | Source         |
|---------------|---------|---------|----------------|
| gf            | 4.18    | kJ/mol  | Joback Method  |
| hf            | -173.31 | kJ/mol  | Joback Method  |
| hfus          | 12.85   | kJ/mol  | Joback Method  |
| hvap          | 46.83   | kJ/mol  | Joback Method  |
| log10ws       | -2.65   |         | Crippen Method |
| logp          | 2.488   |         | Crippen Method |
| mcvol         | 133.870 | ml/mol  | McGowan Method |
| pc            | 3032.27 | kPa     | Joback Method  |
| ripol         | 1653.00 |         | NIST Webbook   |
| tb            | 508.92  | K       | Joback Method  |
| tc            | 722.30  | K       | Joback Method  |
| tf            | 267.64  | K       | Joback Method  |
| vc            | 0.512   | m3/kmol | Joback Method  |

## Temperature Dependent Properties

| Property code | Value     | Unit    | Temperature [K] | Source        |
|---------------|-----------|---------|-----------------|---------------|
| cpg           | 295.28    | J/molxK | 508.92          | Joback Method |
| cpg           | 310.10    | J/molxK | 544.48          | Joback Method |
| cpg           | 324.10    | J/molxK | 580.05          | Joback Method |
| cpg           | 337.29    | J/molxK | 615.61          | Joback Method |
| cpg           | 349.72    | J/molxK | 651.17          | Joback Method |
| cpg           | 361.39    | J/molxK | 686.74          | Joback Method |
| cpg           | 372.35    | J/molxK | 722.30          | Joback Method |
| dvisc         | 0.0040883 | Paxs    | 267.64          | Joback Method |
| dvisc         | 0.0019089 | Paxs    | 307.85          | Joback Method |

|       |           |      |        |               |
|-------|-----------|------|--------|---------------|
| dvisc | 0.0010628 | Paxs | 348.07 | Joback Method |
| dvisc | 0.0006681 | Paxs | 388.28 | Joback Method |
| dvisc | 0.0004582 | Paxs | 428.49 | Joback Method |
| dvisc | 0.0003352 | Paxs | 468.71 | Joback Method |
| dvisc | 0.0002577 | Paxs | 508.92 | 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=R562415&amp;Units=SI">http://webbook.nist.gov/cgi/cbook.cgi?ID=R562415&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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