

# exo-Fenchyl acetate

**Inchi:** InChI=1S/C12H20O2/c1-8(13)14-10-11(2,3)9-5-6-12(10,4)7-9/h9-10H,5-7H2,1-4H3/t9?,1  
**InchiKey:** JUWUWIGZUVEFQB-SQLBVSGCSA-N  
**Formula:** C12H20O2  
**SMILES:** CC(=O)OC1C2(C)CCC(C2)C1(C)C  
**Mol. weight [g/mol]:** 196.29

## Physical Properties

| Property code | Value   | Unit                 | Source         |
|---------------|---------|----------------------|----------------|
| gf            | -100.76 | kJ/mol               | Joback Method  |
| hf            | -406.57 | kJ/mol               | Joback Method  |
| hfus          | 13.34   | kJ/mol               | Joback Method  |
| hvap          | 48.54   | kJ/mol               | Joback Method  |
| log10ws       | -2.88   |                      | Crippen Method |
| logp          | 2.764   |                      | Crippen Method |
| mcvol         | 165.660 | ml/mol               | McGowan Method |
| pc            | 2462.92 | kPa                  | Joback Method  |
| rinpol        | 1231.00 |                      | NIST Webbook   |
| rinpol        | 1230.00 |                      | NIST Webbook   |
| rinpol        | 1230.00 |                      | NIST Webbook   |
| rinpol        | 1232.00 |                      | NIST Webbook   |
| rinpol        | 1232.00 |                      | NIST Webbook   |
| rinpol        | 1214.00 |                      | NIST Webbook   |
| rinpol        | 1233.00 |                      | NIST Webbook   |
| ripol         | 1480.00 |                      | NIST Webbook   |
| tb            | 559.14  | K                    | Joback Method  |
| tc            | 774.82  | K                    | Joback Method  |
| tf            | 368.84  | K                    | Joback Method  |
| vc            | 0.631   | m <sup>3</sup> /kmol | Joback Method  |

## Temperature Dependent Properties

| Property code | Value  | Unit    | Temperature [K] | Source        |
|---------------|--------|---------|-----------------|---------------|
| cpg           | 437.90 | J/molxK | 559.14          | Joback Method |
| cpg           | 456.54 | J/molxK | 595.09          | Joback Method |

|     |        |         |        |               |
|-----|--------|---------|--------|---------------|
| cpg | 474.00 | J/mol×K | 631.03 | Joback Method |
| cpg | 490.47 | J/mol×K | 666.98 | Joback Method |
| cpg | 506.18 | J/mol×K | 702.93 | Joback Method |
| cpg | 521.34 | J/mol×K | 738.88 | Joback Method |
| cpg | 536.18 | J/mol×K | 774.82 | Joback Method |

## Sources

|                        |   |
|------------------------|---|
| <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>                     |
| <b>NIST Webbook:</b>   | <a href="http://webbook.nist.gov/cgi/cbook.cgi?ID=R281522&amp;Units=SI">http://webbook.nist.gov/cgi/cbook.cgi?ID=R281522&amp;Units=SI</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>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>rinpola:</b> | Non-polar retention indices                     |
| <b>ripola:</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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