

# (+)-scapanene

**Inchi:** InChI=1S/C15H24/c1-10-11-6-7-12-13(10)14(2,3)8-5-9-15(11,12)4/h11-13H,1,5-9H2,2-4H  
**InchiKey:** OYBHJKDJMRMSLX-CDKPOMLUSA-N  
**Formula:** C15H24  
**SMILES:** C=C1C2C3CCC1C3(C)CCCC2(C)C  
**Mol. weight [g/mol]:** 204.35

## Physical Properties

| Property code | Value   | Unit                 | Source         |
|---------------|---------|----------------------|----------------|
| gf            | 260.15  | kJ/mol               | Joback Method  |
| hf            | -72.81  | kJ/mol               | Joback Method  |
| hfus          | 13.20   | kJ/mol               | Joback Method  |
| hvap          | 46.31   | kJ/mol               | Joback Method  |
| log10ws       | -4.43   |                      | Crippen Method |
| logp          | 4.415   |                      | Crippen Method |
| mcvol         | 185.330 | ml/mol               | McGowan Method |
| pc            | 2113.89 | kPa                  | Joback Method  |
| rinpol        | 1522.50 |                      | NIST Webbook   |
| rinpol        | 1522.00 |                      | NIST Webbook   |
| ripol         | 1685.00 |                      | NIST Webbook   |
| ripol         | 1664.00 |                      | NIST Webbook   |
| ripol         | 1665.00 |                      | NIST Webbook   |
| ripol         | 1665.00 |                      | NIST Webbook   |
| ripol         | 1664.00 |                      | NIST Webbook   |
| tb            | 561.66  | K                    | Joback Method  |
| tc            | 785.91  | K                    | Joback Method  |
| tf            | 358.59  | K                    | Joback Method  |
| vc            | 0.709   | m <sup>3</sup> /kmol | Joback Method  |

## Temperature Dependent Properties

| Property code | Value  | Unit    | Temperature [K] | Source        |
|---------------|--------|---------|-----------------|---------------|
| cpg           | 501.94 | J/mol×K | 561.66          | Joback Method |
| cpg           | 525.20 | J/mol×K | 599.03          | Joback Method |
| cpg           | 546.85 | J/mol×K | 636.41          | Joback Method |

|     |        |         |        |               |
|-----|--------|---------|--------|---------------|
| cpg | 567.18 | J/mol×K | 673.78 | Joback Method |
| cpg | 586.48 | J/mol×K | 711.16 | Joback Method |
| cpg | 605.02 | J/mol×K | 748.53 | Joback Method |
| cpg | 623.10 | J/mol×K | 785.91 | Joback Method |

## Sources

|                        |   |
|------------------------|---|
| <b>NIST Webbook:</b>   | <a href="http://webbook.nist.gov/cgi/cbook.cgi?ID=R141850&amp;Units=SI">http://webbook.nist.gov/cgi/cbook.cgi?ID=R141850&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>                                     |
| <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>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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