

# (+)-N-Methylephedrine

|                      |                                                                                                                                                             |
|----------------------|-------------------------------------------------------------------------------------------------------------------------------------------------------------|
| Other names:         | (1S,2R)-(+)-N-Methylephedrine<br>Benzenemethanol, «alpha»-[1-(dimethylamino)ethyl]-, [S-(R*,S*)]-<br>L-(+)-erythro-N-methylephedrine<br>d-N-methylephedrine |
| Inchi:               | InChI=1S/C11H17NO/c1-9(12(2)3)11(13)10-7-5-4-6-8-10/h4-9,11,13H,1-3H3                                                                                       |
| InchiKey:            | FMCGSUUBYTWNDP-UHFFFAOYSA-N                                                                                                                                 |
| Formula:             | C11H17NO                                                                                                                                                    |
| SMILES:              | CC(C(O)c1ccccc1)N(C)C                                                                                                                                       |
| Mol. weight [g/mol]: | 179.26                                                                                                                                                      |
| CAS:                 | 42151-56-4                                                                                                                                                  |

## Physical Properties

| Property code | Value   | Unit    | Source         |
|---------------|---------|---------|----------------|
| gf            | 123.23  | kJ/mol  | Joback Method  |
| hf            | -129.10 | kJ/mol  | Joback Method  |
| hfus          | 18.35   | kJ/mol  | Joback Method  |
| hvap          | 60.30   | kJ/mol  | Joback Method  |
| log10ws       | -1.93   |         | Crippen Method |
| logp          | 1.670   |         | Crippen Method |
| mcvol         | 157.940 | ml/mol  | McGowan Method |
| pc            | 2963.34 | kPa     | Joback Method  |
| tb            | 581.50  | K       | Joback Method  |
| tc            | 776.75  | K       | Joback Method  |
| tf            | 303.44  | K       | Joback Method  |
| vc            | 0.569   | m3/kmol | Joback Method  |

## Temperature Dependent Properties

| Property code | Value  | Unit    | Temperature [K] | Source        |
|---------------|--------|---------|-----------------|---------------|
| cpg           | 395.20 | J/mol×K | 581.50          | Joback Method |
| cpg           | 409.59 | J/mol×K | 614.04          | Joback Method |
| cpg           | 423.12 | J/mol×K | 646.58          | Joback Method |
| cpg           | 435.84 | J/mol×K | 679.12          | Joback Method |
| cpg           | 447.78 | J/mol×K | 711.67          | Joback Method |

|     |        |         |        |               |
|-----|--------|---------|--------|---------------|
| cpg | 458.99 | J/mol×K | 744.21 | Joback Method |
| cpg | 469.51 | J/mol×K | 776.75 | Joback Method |

## Sources

|                                                                                                                                                                                                                            |                                                                                                                                               |
|----------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------|-----------------------------------------------------------------------------------------------------------------------------------------------|
| NIST Webbook:                                                                                                                                                                                                              | <a href="http://webbook.nist.gov/cgi/cbook.cgi?ID=C42151564&amp;Units=SI">http://webbook.nist.gov/cgi/cbook.cgi?ID=C42151564&amp;Units=SI</a> |
| Crippen Method:                                                                                                                                                                                                            | <a href="http://pubs.acs.org/doi/abs/10.1021/ci990307l">http://pubs.acs.org/doi/abs/10.1021/ci990307l</a>                                     |
| Crippen Method:                                                                                                                                                                                                            | <a href="https://www.chemeo.com/doc/models/crippen_log10ws">https://www.chemeo.com/doc/models/crippen_log10ws</a>                             |
| Solid-Liquid Equilibria of N-Methylephedrine Enantiomers and Their Mixtures with Two Chiral Ionic N-Methylephedrine Enantiomers and Their Mixtures in Three Chiral Solvents Distinguished by Chain Length: McGowan Method: | <a href="https://www.doi.org/10.1021/acs.jced.8b01245">https://www.doi.org/10.1021/acs.jced.8b01245</a>                                       |
|                                                                                                                                                                                                                            | <a href="https://www.doi.org/10.1021/je1007839">https://www.doi.org/10.1021/je1007839</a>                                                     |
|                                                                                                                                                                                                                            | <a href="https://en.wikipedia.org/wiki/Joback_method">https://en.wikipedia.org/wiki/Joback_method</a>                                         |
|                                                                                                                                                                                                                            | <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>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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