

# Methane, nitroso-

|                      |                             |
|----------------------|-----------------------------|
| Other names:         | Nitrosomethane              |
|                      | CH3NO                       |
| Inchi:               | InChI=1S/CH3NO/c1-2-3/h1H3  |
| InchiKey:            | IMHRONYAKYWGCC-UHFFFAOYSA-N |
| Formula:             | CH3NO                       |
| SMILES:              | CN=O                        |
| Mol. weight [g/mol]: | 45.04                       |
| CAS:                 | 865-40-7                    |

## Physical Properties

| Property code | Value        | Unit   | Source         |
|---------------|--------------|--------|----------------|
| hf            | -232.16      | kJ/mol | Joback Method  |
| hvap          | 26.92        | kJ/mol | Joback Method  |
| ie            | 9.68 ± 0.05  | eV     | NIST Webbook   |
| ie            | 9.80         | eV     | NIST Webbook   |
| ie            | 9.80         | eV     | NIST Webbook   |
| ie            | 9.76 ± 0.05  | eV     | NIST Webbook   |
| ie            | 9.68 ± 0.05  | eV     | NIST Webbook   |
| ie            | 9.30         | eV     | NIST Webbook   |
| ie            | 8.70 ± 0.10  | eV     | NIST Webbook   |
| ie            | 10.80 ± 0.30 | eV     | NIST Webbook   |
| ie            | 9.30         | eV     | NIST Webbook   |
| log10ws       | -0.43        |        | Crippen Method |
| logp          | 0.383        |        | Crippen Method |
| mcvol         | 36.500       | ml/mol | McGowan Method |
| pc            | 5636.27      | kPa    | Joback Method  |
| tb            | 285.68       | K      | Joback Method  |
| tc            | 451.23       | K      | Joback Method  |

## Sources

|                 |                                                                                                                       |
|-----------------|-----------------------------------------------------------------------------------------------------------------------|
| Crippen Method: | <a href="https://www.chemeo.com/doc/models/crippen_log10ws">https://www.chemeo.com/doc/models/crippen_log10ws</a>     |
| Joback Method:  | <a href="https://en.wikipedia.org/wiki/Joback_method">https://en.wikipedia.org/wiki/Joback_method</a>                 |
| McGowan Method: | <a href="http://link.springer.com/article/10.1007/BF02311772">http://link.springer.com/article/10.1007/BF02311772</a> |

**NIST Webbook:**

<http://webbook.nist.gov/cgi/cbook.cgi?ID=C865407&Units=SI>

**Crippen Method:**

<http://pubs.acs.org/doi/abs/10.1021/ci990307l>

## Legend

|                            |                                                 |
|----------------------------|-------------------------------------------------|
| <b>hf:</b>                 | Enthalpy of formation at standard conditions    |
| <b>h<sub>vap</sub>:</b>    | Enthalpy of vaporization at standard conditions |
| <b>ie:</b>                 | Ionization energy                               |
| <b>log<sub>10</sub>ws:</b> | Log10 of Water solubility in mol/l              |
| <b>log<sub>p</sub>:</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                            |

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<https://www.chemeo.com/cid/23-818-8/Methane-nitroso.pdf>

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