

# 1-tert-Butyl-4-nitrobenzene

|                      |                                                                                                                                                                                |
|----------------------|--------------------------------------------------------------------------------------------------------------------------------------------------------------------------------|
| Other names:         | 1-Tert-butyl-2-nitrobenzene<br>4-t-Butylnitrobenzene<br>4-tert-Butylnitrobenzene<br>Benzene, 1-(1,1-dimethylethyl)-4-nitro-<br>p-Nitro-t-butylbenzene<br>p-t-Butylnitrobenzene |
| Inchi:               | InChI=1S/C10H13NO2/c1-10(2,3)8-4-6-9(7-5-8)11(12)13/h4-7H,1-3H3                                                                                                                |
| InchiKey:            | XSCPVQNNFLHGHY-UHFFFAOYSA-N                                                                                                                                                    |
| Formula:             | C10H13NO2                                                                                                                                                                      |
| SMILES:              | CC(C)(C)c1ccc([N+](=O)[O-])cc1                                                                                                                                                 |
| Mol. weight [g/mol]: | 179.22                                                                                                                                                                         |
| CAS:                 | 3282-56-2                                                                                                                                                                      |

## Physical Properties

| Property code | Value       | Unit    | Source         |
|---------------|-------------|---------|----------------|
| gf            | 174.49      | kJ/mol  | Joback Method  |
| hf            | -44.18      | kJ/mol  | Joback Method  |
| hfus          | 19.26       | kJ/mol  | Joback Method  |
| hvap          | 56.09       | kJ/mol  | Joback Method  |
| ie            | 9.29 ± 0.02 | eV      | NIST Webbook   |
| log10ws       | -3.44       |         | Crippen Method |
| logp          | 2.892       |         | Crippen Method |
| mcvol         | 145.420     | ml/mol  | McGowan Method |
| pc            | 2982.79     | kPa     | Joback Method  |
| tb            | 608.47      | K       | Joback Method  |
| tc            | 858.26      | K       | Joback Method  |
| tf            | 387.43      | K       | Joback Method  |
| vc            | 0.558       | m3/kmol | Joback Method  |

## Temperature Dependent Properties

| Property code | Value  | Unit    | Temperature [K] | Source        |
|---------------|--------|---------|-----------------|---------------|
| cpg           | 361.51 | J/mol×K | 608.47          | Joback Method |
| cpg           | 376.25 | J/mol×K | 650.10          | Joback Method |

|     |        |         |        |               |
|-----|--------|---------|--------|---------------|
| cpg | 389.79 | J/mol×K | 691.73 | Joback Method |
| cpg | 402.22 | J/mol×K | 733.36 | Joback Method |
| cpg | 413.64 | J/mol×K | 775.00 | Joback Method |
| cpg | 424.12 | J/mol×K | 816.63 | Joback Method |
| cpg | 433.75 | J/mol×K | 858.26 | Joback Method |

## Pressure Dependent Properties

| Property code | Value         | Unit | Pressure [kPa] | Source       |
|---------------|---------------|------|----------------|--------------|
| tbrp          | 416.00 ± 1.00 | K    | 2.00           | NIST Webbook |

## Correlations

| Information                 | Value                         |
|-----------------------------|-------------------------------|
| Property code               | pvap                          |
| Equation                    | $\ln(P_{vp}) = A + B/(T + C)$ |
| Coeff. A                    | 1.19846e+01                   |
| Coeff. B                    | -3.48795e+03                  |
| Coeff. C                    | -8.13960e+01                  |
| Temperature range (K), min. | 379.59                        |
| Temperature range (K), max. | 604.08                        |

## Sources

|                                      |                                                                                                                                                                                         |
|--------------------------------------|-----------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------|
| 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:                        | <a href="http://webbook.nist.gov/cgi/cbook.cgi?ID=C3282562&amp;Units=SI">http://webbook.nist.gov/cgi/cbook.cgi?ID=C3282562&amp;Units=SI</a>                                             |
| The Yaws Handbook of Vapor Pressure: | <a href="https://www.sciencedirect.com/book/9780128029992/the-yaws-handbook-of-vapor-pressure">https://www.sciencedirect.com/book/9780128029992/the-yaws-handbook-of-vapor-pressure</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>                                                                       |

# 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>ie:</b>      | Ionization energy                               |
| <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>pvap:</b>    | Vapor pressure                                  |
| <b>tb:</b>      | Normal Boiling Point Temperature                |
| <b>tbrp:</b>    | Boiling point at reduced pressure               |
| <b>tc:</b>      | Critical Temperature                            |
| <b>tf:</b>      | Normal melting (fusion) point                   |
| <b>vc:</b>      | Critical Volume                                 |

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