

Benzene, 2-bromo-1-methyl-4-nitro-

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| Other names: | 1-Bromo-2-methyl-5-nitrobenzene 2-Bromo-4-nitrotoluene Toluene, 2-bromo-4-nitro- |
| Inchi: | InChI=1S/C7H6BrNO2/c1-5-2-3-6(9(10)11)4-7(5)8/h2-4H,1H3 |
| InchiKey: | XFZFJQHXXWJIBQV-UHFFFAOYSA-N |
| Formula: | C7H6BrNO2 |
| SMILES: | Cc1ccc([N+](=O)[O-])cc1Br |
| Mol. weight [g/mol]: | 216.03 |
| CAS: | 7745-93-9 |

Physical Properties

| Property code | Value | Unit | Source |
|---------------|---------|---------|----------------|
| gf | 151.08 | kJ/mol | Joback Method |
| hf | 41.35 | kJ/mol | Joback Method |
| hfus | 23.79 | kJ/mol | Joback Method |
| hvap | 57.80 | kJ/mol | Joback Method |
| log10ws | -3.76 | | Crippen Method |
| logp | 2.666 | | Crippen Method |
| mcvol | 120.650 | ml/mol | McGowan Method |
| pc | 4397.41 | kPa | Joback Method |
| tb | 614.20 | K | Joback Method |
| tc | 879.26 | K | Joback Method |
| tf | 423.52 | K | Joback Method |
| vc | 0.464 | m3/kmol | Joback Method |

Temperature Dependent Properties

| Property code | Value | Unit | Temperature [K] | Source |
|---------------|--------|---------|-----------------|---------------|
| cpg | 249.82 | J/molxK | 614.20 | Joback Method |
| cpg | 259.45 | J/molxK | 658.38 | Joback Method |
| cpg | 268.29 | J/molxK | 702.55 | Joback Method |
| cpg | 276.39 | J/molxK | 746.73 | Joback Method |
| cpg | 283.80 | J/molxK | 790.91 | Joback Method |
| cpg | 290.58 | J/molxK | 835.09 | Joback Method |

cpg

296.77

J/mol×K

879.26

Joback Method

Pressure Dependent Properties

| Property code | Value | Unit | Pressure [kPa] | Source |
|---------------|--------|------|----------------|--------------|
| tbrp | 423.70 | K | 2.70 | NIST Webbook |

Sources

| | |
|------------------------|---|
| NIST Webbook: | http://webbook.nist.gov/cgi/cbook.cgi?ID=C7745939&Units=SI |
| Crippen Method: | http://pubs.acs.org/doi/abs/10.1021/ci990307l |
| Crippen Method: | https://www.chemeo.com/doc/models/crippen_log10ws |
| Joback Method: | https://en.wikipedia.org/wiki/Joback_method |
| McGowan Method: | http://link.springer.com/article/10.1007/BF02311772 |

Legend

| | |
|-----------------|---|
| cpg: | Ideal gas heat capacity |
| gf: | Standard Gibbs free energy of formation |
| hf: | Enthalpy of formation at standard conditions |
| hfus: | Enthalpy of fusion at standard conditions |
| hvap: | Enthalpy of vaporization at standard conditions |
| log10ws: | Log10 of Water solubility in mol/l |
| logp: | Octanol/Water partition coefficient |
| mcvol: | McGowan's characteristic volume |
| pc: | Critical Pressure |
| tb: | Normal Boiling Point Temperature |
| tbrp: | Boiling point at reduced pressure |
| tc: | Critical Temperature |
| tf: | Normal melting (fusion) point |
| vc: | Critical Volume |

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<https://www.chemeo.com/cid/19-102-7/Benzene-2-bromo-1-methyl-4-nitro.pdf>

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