

Benzoic acid, 4-methoxy-3-nitro-

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| Other names: | 4-Methoxy-3-nitrobenzoic acid 3-Nitro-4-methoxybenzoic acid p-Anisic acid, 3-nitro- 3-Nitro-p-anisic acid |
| Inchi: | InChI=1S/C8H7NO5/c1-14-7-3-2-5(8(10)11)4-6(7)9(12)13/h2-4H,1H3,(H,10,11) |
| InchiKey: | ANXBDAFDZSXOPQ-UHFFFAOYSA-N |
| Formula: | C8H7NO5 |
| SMILES: | COc1ccc(C(=O)O)cc1[N+](=O)[O-] |
| Mol. weight [g/mol]: | 197.14 |
| CAS: | 89-41-8 |

Physical Properties

| Property code | Value | Unit | Source |
|---------------|---------------|----------------------|----------------|
| gf | -225.56 | kJ/mol | Joback Method |
| hf | -402.65 | kJ/mol | Joback Method |
| hfus | 27.98 | kJ/mol | Joback Method |
| hsub | 131.20 ± 0.80 | kJ/mol | NIST Webbook |
| hvap | 79.43 | kJ/mol | Joback Method |
| log10ws | -2.30 | | Crippen Method |
| logp | 1.302 | | Crippen Method |
| mcpvol | 130.550 | ml/mol | McGowan Method |
| pc | 4244.08 | kPa | Joback Method |
| tb | 739.39 | K | Joback Method |
| tc | 966.92 | K | Joback Method |
| tf | 507.97 | K | Joback Method |
| vc | 0.500 | m ³ /kmol | Joback Method |

Temperature Dependent Properties

| Property code | Value | Unit | Temperature [K] | Source |
|---------------|--------|---------|-----------------|---------------|
| cpg | 336.61 | J/mol×K | 739.39 | Joback Method |
| cpg | 344.79 | J/mol×K | 777.31 | Joback Method |
| cpg | 352.32 | J/mol×K | 815.23 | Joback Method |
| cpg | 359.20 | J/mol×K | 853.16 | Joback Method |

| | | | | |
|-------|---------------|---------|--------|---------------|
| cpg | 365.43 | J/mol×K | 891.08 | Joback Method |
| cpg | 371.04 | J/mol×K | 929.00 | Joback Method |
| cpg | 376.02 | J/mol×K | 966.92 | Joback Method |
| hsubt | 126.50 ± 0.80 | kJ/mol | 394.00 | NIST Webbook |

Sources

| | |
|------------------------|---|
| McGowan Method: | http://link.springer.com/article/10.1007/BF02311772 |
| NIST Webbook: | http://webbook.nist.gov/cgi/cbook.cgi?ID=C89418&Units=SI |
| Crippen Method: | http://pubs.acs.org/doi/abs/10.1021/ci9903071 |
| Crippen Method: | https://www.chemeo.com/doc/models/crippen_log10ws |
| Joback Method: | https://en.wikipedia.org/wiki/Joback_method |

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 |
| hsub: | Enthalpy of sublimation at standard conditions |
| hsubt: | Enthalpy of sublimation at a given temperature |
| hvap: | Enthalpy of vaporization at standard conditions |
| log10ws: | Log10 of Water solubility in mol/l |
| logp: | Octanol/Water partition coefficient |
| mconvol: | McGowan's characteristic volume |
| pc: | Critical Pressure |
| tb: | Normal Boiling Point Temperature |
| tc: | Critical Temperature |
| tf: | Normal melting (fusion) point |
| vc: | Critical Volume |

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