

1,3-Benzenediacetonitrile

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|-----------------------------|---|
| Other names: | m-Benzenediacetonitrile m-Bis(cyanomethyl)benzene m-Phenylenediacetonitrile 1,3-Phenylenediacetonitrile 1,3-Bis(cyanomethyl)benzene m-Xylylene dicyanide Acetonitrile, m-phenylenedi- (3-Cyanomethyl-phenyl)-acetonitrile NSC 77095 |
| Inchi: | InChI=1S/C10H8N2/c11-6-4-9-2-1-3-10(8-9)5-7-12/h1-3,8H,4-5H2 |
| InchiKey: | GRPFZJNUYXIVSL-UHFFFAOYSA-N |
| Formula: | C10H8N2 |
| SMILES: | N#CCc1cccc(CC#N)c1 |
| Mol. weight [g/mol]: | 156.18 |
| CAS: | 626-22-2 |

Physical Properties

| Property code | Value | Unit | Source |
|---------------|---------|----------------------|----------------|
| gf | 402.46 | kJ/mol | Joback Method |
| hf | 305.09 | kJ/mol | Joback Method |
| hfus | 18.32 | kJ/mol | Joback Method |
| hvap | 61.75 | kJ/mol | Joback Method |
| log10ws | -2.81 | | Crippen Method |
| logp | 1.819 | | Crippen Method |
| mcvol | 130.760 | ml/mol | McGowan Method |
| pc | 2735.42 | kPa | Joback Method |
| tb | 664.02 | K | Joback Method |
| tc | 902.09 | K | Joback Method |
| tf | 371.38 | K | Joback Method |
| vc | 0.539 | m ³ /kmol | Joback Method |

Temperature Dependent Properties

| Property code | Value | Unit | Temperature [K] | Source |
|---------------|-------|------|-----------------|--------|
|---------------|-------|------|-----------------|--------|

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|-----|--------|---------|--------|---------------|
| cpg | 304.87 | J/mol×K | 664.02 | Joback Method |
| cpg | 314.46 | J/mol×K | 703.70 | Joback Method |
| cpg | 323.36 | J/mol×K | 743.38 | Joback Method |
| cpg | 331.60 | J/mol×K | 783.05 | Joback Method |
| cpg | 339.22 | J/mol×K | 822.73 | Joback Method |
| cpg | 346.26 | J/mol×K | 862.41 | Joback Method |
| cpg | 352.75 | J/mol×K | 902.09 | Joback Method |

Pressure Dependent Properties

| Property code | Value | Unit | Pressure [kPa] | Source |
|---------------|--------|------|----------------|--------------|
| tbrp | 443.20 | K | 3.30 | NIST Webbook |

Sources

| | |
|------------------------|---|
| 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 |
| NIST Webbook: | http://webbook.nist.gov/cgi/cbook.cgi?ID=C626222&Units=SI |
| Crippen Method: | http://pubs.acs.org/doi/abs/10.1021/ci9903071 |

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 |
| mccvol: | 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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