

# Acetonitrile, (2,6-dichlorophenyl)-

|                      |                                                                                                                                            |
|----------------------|--------------------------------------------------------------------------------------------------------------------------------------------|
| Other names:         | (2,6-Dichlorophenyl)acetonitrile<br>2,6-Dichlorobenzeneacetonitrile<br>2,6-Dichlorobenzyl cyanide<br>2,6-Dichlorophenylacetic acid nitrile |
| Inchi:               | InChI=1S/C8H5Cl2N/c9-7-2-1-3-8(10)6(7)4-5-11/h1-3H,4H2                                                                                     |
| InchiKey:            | AOEJUUCUKRUCF-UHFFFAOYSA-N                                                                                                                 |
| Formula:             | C8H5Cl2N                                                                                                                                   |
| SMILES:              | N#CCc1c(Cl)cccc1Cl                                                                                                                         |
| Mol. weight [g/mol]: | 186.04                                                                                                                                     |
| CAS:                 | 3215-64-3                                                                                                                                  |

## Physical Properties

| Property code | Value       | Unit    | Source         |
|---------------|-------------|---------|----------------|
| gf            | 218.95      | kJ/mol  | Joback Method  |
| hf            | 138.54      | kJ/mol  | Joback Method  |
| hfus          | 19.64       | kJ/mol  | Joback Method  |
| hvap          | 56.25       | kJ/mol  | Joback Method  |
| ie            | 9.73 ± 0.06 | eV      | NIST Webbook   |
| log10ws       | -3.51       |         | Crippen Method |
| logp          | 3.059       |         | Crippen Method |
| mcvol         | 125.680     | ml/mol  | McGowan Method |
| pc            | 3142.03     | kPa     | Joback Method  |
| tb            | 596.02      | K       | Joback Method  |
| tc            | 838.45      | K       | Joback Method  |
| tf            | 356.21      | K       | Joback Method  |
| vc            | 0.499       | m3/kmol | Joback Method  |

## Temperature Dependent Properties

| Property code | Value  | Unit    | Temperature [K] | Source        |
|---------------|--------|---------|-----------------|---------------|
| cpg           | 241.66 | J/mol×K | 596.02          | Joback Method |
| cpg           | 249.88 | J/mol×K | 636.43          | Joback Method |
| cpg           | 257.50 | J/mol×K | 676.83          | Joback Method |
| cpg           | 264.56 | J/mol×K | 717.24          | Joback Method |

|     |        |         |        |               |
|-----|--------|---------|--------|---------------|
| cpg | 271.08 | J/mol×K | 757.64 | Joback Method |
| cpg | 277.09 | J/mol×K | 798.05 | Joback Method |
| cpg | 282.61 | J/mol×K | 838.45 | Joback Method |

## Sources

|                        |                                                                                                                                             |
|------------------------|---------------------------------------------------------------------------------------------------------------------------------------------|
| <b>Crippen Method:</b> | <a href="https://www.chemeo.com/doc/models/crippen_log10ws">https://www.chemeo.com/doc/models/crippen_log10ws</a>                           |
| <b>Joback Method:</b>  | <a href="https://en.wikipedia.org/wiki/Joback_method">https://en.wikipedia.org/wiki/Joback_method</a>                                       |
| <b>McGowan Method:</b> | <a href="http://link.springer.com/article/10.1007/BF02311772">http://link.springer.com/article/10.1007/BF02311772</a>                       |
| <b>NIST Webbook:</b>   | <a href="http://webbook.nist.gov/cgi/cbook.cgi?ID=C3215643&amp;Units=SI">http://webbook.nist.gov/cgi/cbook.cgi?ID=C3215643&amp;Units=SI</a> |
| <b>Crippen Method:</b> | <a href="http://pubs.acs.org/doi/abs/10.1021/ci990307l">http://pubs.acs.org/doi/abs/10.1021/ci990307l</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>tb:</b>      | Normal Boiling Point Temperature                |
| <b>tc:</b>      | Critical Temperature                            |
| <b>tf:</b>      | Normal melting (fusion) point                   |
| <b>vc:</b>      | Critical Volume                                 |

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