

2,5-dimethyl-4-ethylpyridine

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|-----------------------------|---|
| Inchi: | InChI=1S/C9H13N/c1-4-9-5-8(3)10-6-7(9)2/h5-6H,4H2,1-3H3 |
| InchiKey: | JBNJDXYIBJKBL-UHFFFAOYSA-N |
| Formula: | C9H13N |
| SMILES: | CCc1cc(C)ncc1C |
| Mol. weight [g/mol]: | 135.21 |

Physical Properties

| Property code | Value | Unit | Source |
|---------------|---------|--------|----------------|
| log10ws | -2.99 | | Crippen Method |
| logp | 2.261 | | Crippen Method |
| mcvol | 123.890 | ml/mol | McGowan Method |
| rinpol | 1167.00 | | NIST Webbook |
| rinpol | 1167.00 | | NIST Webbook |

Sources

| | |
|------------------------|---|
| Crippen Method: | https://www.chemeo.com/doc/models/crippen_log10ws |
| McGowan Method: | http://link.springer.com/article/10.1007/BF02311772 |
| NIST Webbook: | http://webbook.nist.gov/cgi/cbook.cgi?ID=R142284&Units=SI |
| Crippen Method: | http://pubs.acs.org/doi/abs/10.1021/ci990307l |

Legend

| | |
|-----------------|-------------------------------------|
| log10ws: | Log10 of Water solubility in mol/l |
| logp: | Octanol/Water partition coefficient |
| mcvol: | McGowan's characteristic volume |
| rinpol: | Non-polar retention indices |

Latest version available from:

<https://www.chemeo.com/cid/52-211-9/2-5-dimethyl-4-ethylpyridine.pdf>

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