

# 7H-Benzo[c]carbazole

|                      |                                                                                  |
|----------------------|----------------------------------------------------------------------------------|
| Other names:         | 3,4-Benzocarbazole<br>7-Aza-7H-benzo[c]fluorene                                  |
| Inchi:               | InChI=1S/C16H11N/c1-2-6-12-11(5-1)9-10-15-16(12)13-7-3-4-8-14(13)17-15/h1-10,17H |
| InchiKey:            | UGFOTZLGPPWNPY-UHFFFAOYSA-N                                                      |
| Formula:             | C16H11N                                                                          |
| SMILES:              | c1ccc2c(c1)ccc1[nH]c3ccccc3c12                                                   |
| Mol. weight [g/mol]: | 217.27                                                                           |
| CAS:                 | 205-25-4                                                                         |

## Physical Properties

| Property code | Value   | Unit   | Source         |
|---------------|---------|--------|----------------|
| log10ws       | -6.10   |        | Crippen Method |
| logp          | 3.992   |        | Crippen Method |
| mcvol         | 168.440 | ml/mol | McGowan Method |

## Sources

|                 |                                                                                                                                           |
|-----------------|-------------------------------------------------------------------------------------------------------------------------------------------|
| McGowan Method: | <a href="http://link.springer.com/article/10.1007/BF02311772">http://link.springer.com/article/10.1007/BF02311772</a>                     |
| NIST Webbook:   | <a href="http://webbook.nist.gov/cgi/cbook.cgi?ID=C205254&amp;Units=SI">http://webbook.nist.gov/cgi/cbook.cgi?ID=C205254&amp;Units=SI</a> |
| Crippen Method: | <a href="http://pubs.acs.org/doi/abs/10.1021/ci990307l">http://pubs.acs.org/doi/abs/10.1021/ci990307l</a>                                 |
| Crippen Method: | <a href="https://www.chemeo.com/doc/models/crippen_log10ws">https://www.chemeo.com/doc/models/crippen_log10ws</a>                         |

## Legend

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

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