

# Cytidine

|                      |                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                     |
|----------------------|-------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------|
| Other names:         | 1«beta»-D-Ribofuranosylcytosine<br>1«beta»-Ribofuranosylcytosine<br>1Â«betaÂ»-D-Ribofuranosylcytosine<br>1Â«betaÂ»-Ribofuranosylcytosine<br>2(1H)-Pyrimidinone, 4-amino-1-«beta»-D-ribofuranosyl-<br>2(1H)-Pyrimidinone, 4-amino-1-Â«betaÂ»-D-ribofuranosyl-<br>4-Amino-1«beta»-D-ribofuranosyl-2(1H)-pyrimidinone<br>4-Amino-1Â«betaÂ»-D-ribofuranosyl-2(1H)-pyrimidinone<br>Cyd<br>Cytosine riboside<br>Cytosine, 1-«beta»-D-ribofuranosyl-<br>Cytosine, 1-Â«betaÂ»-D-ribofuranosyl-<br>NSC 20258<br>«beta»-D-Ribofuranoside, cytosine-1<br>Â«betaÂ»-D-Ribofuranoside, cytosine-1 |
| Inchi:               | InChI=1S/C9H13N3O5/c10-5-1-2-12(9(16)11-5)8-7(15)6(14)4(3-13)17-8/h1-2,4,6-8,13-15                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                  |
| InchiKey:            | UHDGCWIWMRVCDJ-XVFCMESISA-N                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                         |
| Formula:             | C9H13N3O5                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                           |
| SMILES:              | Nc1ccn(C2OC(CO)C(O)C2O)c(=O)n1                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                      |
| Mol. weight [g/mol]: | 243.22                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                              |
| CAS:                 | 65-46-3                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                             |

## Physical Properties

| Property code | Value   | Unit   | Source         |
|---------------|---------|--------|----------------|
| affp          | 982.50  | kJ/mol | NIST Webbook   |
| basg          | 950.00  | kJ/mol | NIST Webbook   |
| log10ws       | 0.62    |        | Crippen Method |
| logp          | -2.563  |        | Crippen Method |
| mcvol         | 162.340 | ml/mol | McGowan Method |

## Temperature Dependent Properties

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

|      |         |       |        |                                                                                                                                                                        |
|------|---------|-------|--------|------------------------------------------------------------------------------------------------------------------------------------------------------------------------|
| rhos | 1530.00 | kg/m3 | 298.15 | Saturation molalities and standard molar enthalpies of solution of cytidine(cr), hypoxanthine(cr), thymidine(cr), thymine(cr), uridine(cr), and xanthine(cr) in H2O(l) |
|------|---------|-------|--------|------------------------------------------------------------------------------------------------------------------------------------------------------------------------|

## Sources

The partial molar volumes at T = (288.15 to 313.15) K, and the partial molar isobaric expansion coefficients for some nucleosides in aqueous solution and adenosine in aqueous solution:

Volumetric properties of the nucleosides adenosine, cytidine, and uridine in aqueous solution. Solution behavior of some nucleic acid bases and nucleosides in water and in aqueous guanidine hydrochloride solutions: Viscometric, calorimetric and spectroscopic compressions of the Nucleosides Adenosine, Cytidine, and Uridine in Aqueous Solution at 298.15 K: Volumetric studies on nucleic acid bases and nucleosides in aqueous Guanidine Hydrochloride solutions at T = (288.15 to 318.15) K and at atmospheric pressure

<https://www.doi.org/10.1016/j.jct.2008.01.020>  
<https://www.doi.org/10.1016/j.jct.2012.12.014>  
<http://link.springer.com/article/10.1007/BF02311772>  
<https://www.doi.org/10.1016/j.jct.2012.12.028>  
<https://www.doi.org/10.1016/j.jct.2015.11.029>  
<http://pubs.acs.org/doi/abs/10.1021/ci990307l>  
<https://www.doi.org/10.1021/je101264r>  
<http://webbook.nist.gov/cgi/cbook.cgi?ID=C65463&Units=SI>  
<https://www.doi.org/10.1016/j.jct.2014.10.015>  
[https://www.chemeo.com/doc/models/crippen\\_log10ws](https://www.chemeo.com/doc/models/crippen_log10ws)  
<https://www.doi.org/10.1016/j.jct.2004.04.005>

## Legend

- affp: Proton affinity
- basg: Gas basicity
- log10ws: Log10 of Water solubility in mol/l
- logp: Octanol/Water partition coefficient
- mcvol: McGowan's characteristic volume
- rhos: Solid Density

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<https://www.chemeo.com/cid/27-883-2/Cytidine.pdf>

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