

Uridine

Other names:

1-((2R,3R,4S,5R)-3,4-dihydroxy-5-(hydroxymethyl)tetrahydrofuran-2-yl)pyrimidine-2,4(1H,3H)-pyrimidinedione
1-.beta.-D-ribofuranosyl-2,4(1H,3H)-pyrimidinedione
1-«beta»-D-Ribofuranosyl-2,4(1H,3H)-pyrimidinedione
1-«beta»-D-Ribofuranosyluracil
NSC 20256
Uracil riboside
Uracil, 1-«beta»-D-ribofuranosyl-
Uracil-1-«beta»-d-ribofuranoside
Urd
Uridin
d-Ribosyl uracil
«beta»-D-Ribofuranoside, 2,4(1H,3H)-pyrimidinedione-1
«beta»-Uridine

Inchi: InChI=1S/C9H12N2O6/c12-3-4-6(14)7(15)8(17-4)11-2-1-5(13)10-9(11)16/h1-2,4,6-8,12,14-16
InchiKey: DRTQHJPMGBCUFC-XVFCMESISA-N
Formula: C9H12N2O6
SMILES: O=c1ccn(C2OC(CO)C(O)C2O)c(=O)[nH]1
Mol. weight [g/mol]: 244.20
CAS: 58-96-8

Physical Properties

Property code	Value	Unit	Source
affp	947.60	kJ/mol	NIST Webbook
affp	899.00	kJ/mol	NIST Webbook
basg	916.60	kJ/mol	NIST Webbook
log10ws	1.38		Crippen Method
logp	-3.334		Crippen Method
mcpol	158.230	ml/mol	McGowan Method

Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
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cps	292.34	J/mol×K	298.15	Molar Heat Capacities of Some Derivatives of Uridine and 2'-Deoxyuridine
cps	294.26	J/mol×K	303.15	Molar Heat Capacities of Some Derivatives of Uridine and 2'-Deoxyuridine
cps	296.18	J/mol×K	308.15	Molar Heat Capacities of Some Derivatives of Uridine and 2'-Deoxyuridine
cps	298.10	J/mol×K	313.15	Molar Heat Capacities of Some Derivatives of Uridine and 2'-Deoxyuridine
cps	300.02	J/mol×K	318.15	Molar Heat Capacities of Some Derivatives of Uridine and 2'-Deoxyuridine
cps	301.93	J/mol×K	323.15	Molar Heat Capacities of Some Derivatives of Uridine and 2'-Deoxyuridine
cps	303.86	J/mol×K	328.15	Molar Heat Capacities of Some Derivatives of Uridine and 2'-Deoxyuridine
cps	305.77	J/mol×K	333.15	Molar Heat Capacities of Some Derivatives of Uridine and 2'-Deoxyuridine
cps	307.69	J/mol×K	338.15	Molar Heat Capacities of Some Derivatives of Uridine and 2'-Deoxyuridine
cps	309.61	J/mol×K	343.15	Molar Heat Capacities of Some Derivatives of Uridine and 2'-Deoxyuridine
cps	311.53	J/mol×K	348.15	Molar Heat Capacities of Some Derivatives of Uridine and 2'-Deoxyuridine
cps	313.45	J/mol×K	353.15	Molar Heat Capacities of Some Derivatives of Uridine and 2'-Deoxyuridine

cps	315.37	J/molxK	358.15	Molar Heat Capacities of Some Derivatives of Uridine and 2'-Deoxyuridine
cps	317.29	J/molxK	363.15	Molar Heat Capacities of Some Derivatives of Uridine and 2'-Deoxyuridine
cps	319.21	J/molxK	368.15	Molar Heat Capacities of Some Derivatives of Uridine and 2'-Deoxyuridine
rhos	1600.00	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

Volumetric interaction coefficients for some nucleosides in aqueous solution at T = 298.15 K: Molar Heat Capacities of Some Derivatives of Uridine and 2'-Deoxyuridine:
Crippen Method:

<https://www.doi.org/10.1016/j.jct.2012.12.014>

Crippen Method:

<https://www.doi.org/10.1021/je800243y>

<http://link.springer.com/article/10.1007/BF02311772>

https://www.chemeo.com/doc/models/crippen_log10ws

The partial molar volumes at T = (288.15 to 313.15) K, and the partial molar heat capacities and expansions

<https://www.doi.org/10.1016/j.jct.2008.01.020>

at T = 298.15 K of cytidine, uridine, and adenosine in aqueous solution.

<http://pubs.acs.org/doi/abs/10.1021/ci9903071>

Volumetric studies on nucleic acid bases and nucleosides in aqueous solution.

<https://www.doi.org/10.1016/j.jct.2014.10.015>

Energy characterization of binary and ternary aqueous solutions at T = 298.15 K of cytidine, uridine, and adenosine compounds.

<https://www.doi.org/10.1016/j.jct.2018.04.013>

Partial Molar Isoentropic and Isothermal Compressions of the Nucleosides Cytidine, Uridine, and Adenosine.

<https://www.doi.org/10.1021/je101264r>

Saturation molalities and standard molar enthalpies of solution of cytidine(cr), hypoxanthine(cr), thymidine(cr), thymine(cr), uridine(cr), and xanthine(cr) in aqueous solution at T = 298.15 K and P = (10 to 120) MPa.

<https://www.doi.org/10.1016/j.jct.2004.04.005>

Solvation behavior of some nucleic acid bases and nucleosides in water and in aqueous guanidine hydrochloride solutions: Viscometric, calorimetric and spectroscopic approach.

<https://www.doi.org/10.1016/j.jct.2012.12.028>

<http://webbook.nist.gov/cgi/cbook.cgi?ID=C58968&Units=SI>

<https://www.doi.org/10.1016/j.jct.2015.11.029>

Legend

affp:	Proton affinity
basg:	Gas basicity
cps:	Solid phase heat capacity
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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