ACYCLOVIR

Other names: 2-Amino-1,9-dihydro-9-((2-hydroxyethoxy)methyl)-6H-purin-6-one
2-Amino-9-[(2-hydroxyethoxy)methyl]-1,9-dihydro-6H-purin-6-one
6H-Purin-6-one, 1,9-dihydro-2-amino-9-((2-hydroxyethoxy)methyl)-
6H-Purin-6-one, 2-amino-1,9-dihydro-9-[(2-hydroxyethoxy)methyl]-
9-(2-Hydroxyethoxymethyl)guanine
9-[(2-hydroxyethoxyl)methyl]guanine
ACV
Acicloftal
Aciclovir
Acyclo-V
Acycloguanosine
BW 248U
Cargosil
NSC 645011
Poviral
Vipral
Virorax
W-248-U
Wellcome-248U
Zovirax
Zyclir

Inchi: InChI=1S/C8H11N5O3/c9-8-11-6-5(7(15)12-8)10-3-13(6)4-16-2-1-14/h3,14H,1-2,4H2,(H

InchiKey: MKUXAQIIEYXACX-UHFFFAOYSA-N

Formula: C8H11N5O3

SMILES: Nc1nc2c(ncn2COCCO)c(=O)[nH]1

Mol. weight [g/mol]: 225.21

Physical Properties

Property code	Value	Unit	Source
log10ws	-1.61		Cheméo Relay (1.0)
logp	-1.814		Crippen Method
mcvol	152.170	ml/mol	McGowan Method
tf	512.83	K	Cheméo Relay (1.0)

Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
hfust	30.44	kJ/mol	528.20	NIST Webbook

Sources

Thermal and spectral characterization of a binary mixture (acyclovir and fluocinolone acetonide): Eutectic reaction and inclusion complexes with beta-cyclodextrin:	https://www.doi.org/10.1016/j.tca.2013.03.013
NIST Webbook:	http://webbook.nist.gov/cgi/cbook.cgi?ID=C59277893&Units=SI
McGowan Method:	http://link.springer.com/article/10.1007/BF02311772
Crippen Method:	http://pubs.acs.org/doi/abs/10.1021/ci990307l
Crippen Method:	https://www.chemeo.com/doc/models/crippen_log10ws
Cheméo Relay (1.0):	
Cheméo Relay (1.0, beta):	

Legend

hfust:	Enthalpy of fusion at a given temperature
log10ws:	Log10 of Water solubility in mol/l
logp:	Octanol/Water partition coefficient
mcvol:	McGowan's characteristic volume
tf:	Normal melting (fusion) point