

# 5«beta»-Pregn-8-ene-3«alpha»,20«alpha»-diol

**Inchi:** InChI=1S/C21H34O2/c1-13(22)17-6-7-18-16-5-4-14-12-15(23)8-10-20(14,2)19(16)9-11-2  
**InchiKey:** CVXRELSVEQTOQH-VDDUQTHDSA-N  
**Formula:** C21H34O2  
**SMILES:** CC(O)C1CCC2C3=C(CCC21C)C1(C)CCC(O)CC1CC3  
**Mol. weight [g/mol]:** 318.49

## Physical Properties

Property code	Value	Unit	Source
gf	16.66	kJ/mol	Joback Method
hf	-501.47	kJ/mol	Joback Method
hfus	26.83	kJ/mol	Joback Method
hvap	94.52	kJ/mol	Joback Method
log10ws	-5.59		Crippen Method
logp	4.451		Crippen Method
mcvol	270.750	ml/mol	McGowan Method
pc	1784.86	kPa	Joback Method
rinpol	2741.00		NIST Webbook
rinpol	2741.00		NIST Webbook
tb	912.37	K	Joback Method
tc	1130.79	K	Joback Method
tf	552.35	K	Joback Method
vc	1.012	m <sup>3</sup> /kmol	Joback Method

## Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	999.40	J/mol×K	912.37	Joback Method
cpg	1023.77	J/mol×K	948.77	Joback Method
cpg	1048.69	J/mol×K	985.18	Joback Method
cpg	1074.44	J/mol×K	1021.58	Joback Method
cpg	1101.32	J/mol×K	1057.98	Joback Method
cpg	1129.61	J/mol×K	1094.38	Joback Method
cpg	1159.60	J/mol×K	1130.79	Joback Method

# Sources

<b>Crippen Method:</b>	<a href="http://pubs.acs.org/doi/abs/10.1021/ci9903071">http://pubs.acs.org/doi/abs/10.1021/ci9903071</a>
<b>Crippen Method:</b>	<a href="https://www.cheméo.com/doc/models/crippen_log10ws">https://www.cheméo.com/doc/models/crippen_log10ws</a>
<b>Joback Method:</b>	<a href="https://en.wikipedia.org/wiki/Joback_method">https://en.wikipedia.org/wiki/Joback_method</a>
<b>McGowan Method:</b>	<a href="http://link.springer.com/article/10.1007/BF02311772">http://link.springer.com/article/10.1007/BF02311772</a>
<b>NIST Webbook:</b>	<a href="http://webbook.nist.gov/cgi/cbook.cgi?ID=R304051&amp;Units=SI">http://webbook.nist.gov/cgi/cbook.cgi?ID=R304051&amp;Units=SI</a>

# Legend

<b>cpg:</b>	Ideal gas heat capacity
<b>gf:</b>	Standard Gibbs free energy of formation
<b>hf:</b>	Enthalpy of formation at standard conditions
<b>hfus:</b>	Enthalpy of fusion at standard conditions
<b>hvp:</b>	Enthalpy of vaporization at standard conditions
<b>log10ws:</b>	Log10 of Water solubility in mol/l
<b>logp:</b>	Octanol/Water partition coefficient
<b>mcvol:</b>	McGowan's characteristic volume
<b>pc:</b>	Critical Pressure
<b>rinp:</b>	Non-polar retention indices
<b>tb:</b>	Normal Boiling Point Temperature
<b>tc:</b>	Critical Temperature
<b>tf:</b>	Normal melting (fusion) point
<b>vc:</b>	Critical Volume

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