

# Acenaphthylene, 1,2-dihydro-5-pentadecyl-

<b>Other names:</b>	5-Pentadecylacenaphthene
<b>Inchi:</b>	InChI=1S/C27H40/c1-2-3-4-5-6-7-8-9-10-11-12-13-14-16-23-19-20-25-22-21-24-17-15-1
<b>InchiKey:</b>	BTWSHSRQUSIAIC-UHFFFAOYSA-N
<b>Formula:</b>	C27H40
<b>SMILES:</b>	CCCCCCCCCCCCCCCCc1ccc2c3c(cccc13)CC2
<b>Mol. weight [g/mol]:</b>	364.61
<b>CAS:</b>	55334-13-9

## Physical Properties

Property code	Value	Unit	Source
gf	447.19	kJ/mol	Joback Method
hf	-108.12	kJ/mol	Joback Method
hfus	54.74	kJ/mol	Joback Method
hvap	81.65	kJ/mol	Joback Method
log10ws	-10.18		Crippen Method
logp	8.572		Crippen Method
mcvol	337.210	ml/mol	McGowan Method
pc	1013.59	kPa	Joback Method
tb	884.90	K	Joback Method
tc	1090.15	K	Joback Method
tf	516.43	K	Joback Method
vc	1.327	m3/kmol	Joback Method

## Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	1100.81	J/molxK	884.90	Joback Method
cpg	1120.85	J/molxK	919.11	Joback Method
cpg	1140.07	J/molxK	953.32	Joback Method
cpg	1158.56	J/molxK	987.53	Joback Method
cpg	1176.47	J/molxK	1021.73	Joback Method
cpg	1193.89	J/molxK	1055.94	Joback Method
cpg	1210.96	J/molxK	1090.15	Joback Method
dvisc	0.0013993	Paxs	516.43	Joback Method

dvisc	0.0009361	Paxs	577.84	Joback Method
dvisc	0.0006765	Paxs	639.25	Joback Method
dvisc	0.0005176	Paxs	700.67	Joback Method
dvisc	0.0004134	Paxs	762.08	Joback Method
dvisc	0.0003415	Paxs	823.49	Joback Method
dvisc	0.0002897	Paxs	884.90	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=C55334139&amp;Units=SI">http://webbook.nist.gov/cgi/cbook.cgi?ID=C55334139&amp;Units=SI</a>

## Legend

<b>cpg:</b>	Ideal gas heat capacity
<b>dvisc:</b>	Dynamic viscosity
<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>hvap:</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>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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