

# dopentacontane

**Inchi:** InChI=1S/C52H106/c1-3-5-7-9-11-13-15-17-19-21-23-25-27-29-31-33-35-37-39-41-43-45  
**InchiKey:** HAIPLQZWDQVPPR-UHFFFAOYSA-N  
**Formula:** C52H106  
**SMILES:** CC  
**Mol. weight [g/mol]:** 731.40  
**CAS:** 7719-79-1

## Physical Properties

Property code	Value	Unit	Source
gf	386.96	kJ/mol	Joback Method
hf	-1116.61	kJ/mol	Joback Method
hfus	130.44	kJ/mol	Joback Method
hvap	261.80 ± 1.50	kJ/mol	NIST Webbook
log10ws	-21.59		Crippen Method
logp	20.531		Crippen Method
mcvol	743.540	ml/mol	McGowan Method
pc	259.31	kPa	Joback Method
tb	1389.16	K	Joback Method
tc	2411.65	K	Joback Method
tf	367.00 ± 5.00	K	NIST Webbook
vc	2.947	m3/kmol	Joback Method

## Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	3786.42	J/molxK	2241.24	Joback Method
cpg	4086.78	J/molxK	2411.65	Joback Method
cpg	3039.24	J/molxK	1389.16	Joback Method
cpg	3148.20	J/molxK	1559.58	Joback Method
cpg	3259.35	J/molxK	1729.99	Joback Method
cpg	3390.75	J/molxK	1900.41	Joback Method
cpg	3560.42	J/molxK	2070.82	Joback Method
dvisc	0.0000009	Paxs	1270.27	Joback Method
dvisc	0.0000006	Paxs	1389.16	Joback Method

dvisc	0.0000340	Paxs	675.80	Joback Method
dvisc	0.0000107	Paxs	794.69	Joback Method
dvisc	0.0000046	Paxs	913.59	Joback Method
dvisc	0.0000024	Paxs	1032.48	Joback Method
dvisc	0.0000014	Paxs	1151.37	Joback Method
hfust	171.80	kJ/mol	366.70	NIST Webbook
hvapt	152.00	kJ/mol	736.00	NIST Webbook
hvapt	261.80	kJ/mol	298.15	Hypothetical Thermodynamic Properties: Vapor Pressures and Vaporization Enthalpies of the Even n-Alkanes from C40 to C76 at T = 298.15 K by Correlation-Gas Chromatography. Are the Vaporization Enthalpies a Linear Function of Carbon Number?

## Sources

**NIST Webbook:** <http://webbook.nist.gov/cgi/cbook.cgi?ID=C7719791&Units=SI>

**Crippen Method:** <http://pubs.acs.org/doi/abs/10.1021/ci990307I>

**Crippen Method:** [https://www.chemeo.com/doc/models/crippen\\_log10ws](https://www.chemeo.com/doc/models/crippen_log10ws)

**Hypothetical Thermodynamic Properties: Vapor Pressures and Vaporization Enthalpies of the Even n-Alkanes from C40 to C76 at T = 298.15 K by Correlation-Gas Chromatography. Are the Vaporization Enthalpies a Linear Function of Carbon Number?:** <https://www.doi.org/10.1021/je7005852>

**Hypothetical Thermodynamic Properties: Vapor Pressures and Vaporization Enthalpies of the Even n-Alkanes from C40 to C76 at T = 298.15 K by Correlation-Gas Chromatography. Are the Vaporization Enthalpies a Linear Function of Carbon Number?:** [https://en.wikipedia.org/wiki/Joback\\_method](https://en.wikipedia.org/wiki/Joback_method)

**Hypothetical Thermodynamic Properties: Vapor Pressures and Vaporization Enthalpies of the Even n-Alkanes from C40 to C76 at T = 298.15 K by Correlation-Gas Chromatography. Are the Vaporization Enthalpies a Linear Function of Carbon Number?:** <http://link.springer.com/article/10.1007/BF02311772>

## 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>hfust:</b>	Enthalpy of fusion at a given temperature
<b>hvap:</b>	Enthalpy of vaporization at standard conditions

<b>hvapt:</b>	Enthalpy of vaporization at a given temperature
<b>log10ws:</b>	Log10 of Water solubility in mol/l
<b>logp:</b>	Octanol/Water partition coefficient
<b>mccvol:</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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