

Octacosane

Other names:	n-Octacosane
Inchi:	InChI=1S/C28H58/c1-3-5-7-9-11-13-15-17-19-21-23-25-27-28-26-24-22-20-18-16-14-12
InchiKey:	ZYURHZPYMFLWSH-UHFFFAOYSA-N
Formula:	C28H58
SMILES:	CCCCCCCCCCCCCCCCCCCCCCCCCCCCCC
Mol. weight [g/mol]:	394.76
CAS:	630-02-4

Physical Properties

Property code	Value	Unit	Source
gf	184.88	kJ/mol	Joback Method
hf	-621.25	kJ/mol	Joback Method
hfus	35.53	kJ/mol	Measurement of enthalpy curves of phase change materials via DSC and T-History: When are both methods needed to estimate the behaviour of the bulk material in applications?
hsub	209.00 ± 10.00	kJ/mol	NIST Webbook
hvap	141.90	kJ/mol	NIST Webbook
hvap	152.40 ± 2.90	kJ/mol	NIST Webbook
hvap	150.80 ± 0.50	kJ/mol	NIST Webbook
hvap	150.70 ± 1.70	kJ/mol	NIST Webbook
log10ws	-11.54		Crippen Method
logp	11.169		Crippen Method
mcpol	405.380	ml/mol	McGowan Method
pc	655.44	kPa	Joback Method
rinpol	461.80		NIST Webbook
rinpol	442.42		NIST Webbook
rinpol	461.80		NIST Webbook
tb	704.80	K	NIST Webbook
tc	836.00	K	Critical temperatures and pressures of C40, C44, and C60 normal alkanes measured by the pulse-heating technique
tf	334.55	K	KDB
tt	333.98 ± 0.70	K	NIST Webbook

tt

330.40

K

Thermal conductivity of
heavy, even-carbon
number n-alkanes (C22 to
C32)

vc

1.603

m3/kmol

Joback Method

Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	1347.97	J/mol×K	840.04	Joback Method
cpg	1482.44	J/mol×K	1030.18	Joback Method
cpg	1462.92	J/mol×K	998.49	Joback Method
cpg	1442.33	J/mol×K	966.80	Joback Method
cpg	1420.59	J/mol×K	935.11	Joback Method
cpg	1397.66	J/mol×K	903.42	Joback Method
cpg	1373.48	J/mol×K	871.73	Joback Method
cpl	937.00	J/mol×K	353.00	NIST Webbook
dvisc	0.0003995	Paxs	573.15	Liquid Viscosity and Surface Tension of n-Dodecane, n-Octacosane, Their Mixtures, and a Wax between 323 and 573 K by Surface Light Scattering
dvisc	0.0004753	Paxs	548.15	Liquid Viscosity and Surface Tension of n-Dodecane, n-Octacosane, Their Mixtures, and a Wax between 323 and 573 K by Surface Light Scattering
dvisc	0.0006847	Paxs	498.15	Liquid Viscosity and Surface Tension of n-Dodecane, n-Octacosane, Their Mixtures, and a Wax between 323 and 573 K by Surface Light Scattering

dvisc	0.0008318	Paxs	473.15	Liquid Viscosity and Surface Tension of n-Dodecane, n-Octacosane, Their Mixtures, and a Wax between 323 and 573 K by Surface Light Scattering
dvisc	0.0010252	Paxs	448.15	Liquid Viscosity and Surface Tension of n-Dodecane, n-Octacosane, Their Mixtures, and a Wax between 323 and 573 K by Surface Light Scattering
dvisc	0.0013609	Paxs	423.15	Liquid Viscosity and Surface Tension of n-Dodecane, n-Octacosane, Their Mixtures, and a Wax between 323 and 573 K by Surface Light Scattering
dvisc	0.0018142	Paxs	398.15	Liquid Viscosity and Surface Tension of n-Dodecane, n-Octacosane, Their Mixtures, and a Wax between 323 and 573 K by Surface Light Scattering
dvisc	0.0005675	Paxs	523.15	Liquid Viscosity and Surface Tension of n-Dodecane, n-Octacosane, Their Mixtures, and a Wax between 323 and 573 K by Surface Light Scattering
hfust	67.38	kJ/mol	334.20	NIST Webbook
hfust	63.00	kJ/mol	333.40	NIST Webbook
hfust	64.64	kJ/mol	334.50	NIST Webbook
hfust	35.44	kJ/mol	331.30	NIST Webbook
hfust	64.64	kJ/mol	334.50	NIST Webbook
hsubt	195.80 ± 2.20	kJ/mol	326.00	NIST Webbook
hvapt	106.60	kJ/mol	593.00	NIST Webbook

hvapt	141.90	kJ/mol	298.15	Vapor Pressures and Vaporization Enthalpies of the n-Alkanes from C21 to C30 at T = 298.15 K by Correlation Gas Chromatography
hvapt	117.40 ± 1.20	kJ/mol	375.50	NIST Webbook
hvapt	118.50	kJ/mol	435.50	NIST Webbook
hvapt	100.50	kJ/mol	535.50	NIST Webbook
hvapt	128.60	kJ/mol	298.00	A Comparison of Results by Correlation Gas Chromatography with Another Gas Chromatographic Retention Time Technique. The Effects of Retention Time Coincidence on Vaporization Enthalpy and Vapor Pressure
hvapt	105.50	kJ/mol	459.50	NIST Webbook
hvapt	103.10 ± 3.00	kJ/mol	494.00	NIST Webbook
hvapt	99.27 ± 0.06	kJ/mol	510.00	NIST Webbook
hvapt	100.60	kJ/mol	512.50	NIST Webbook
hvapt	98.10	kJ/mol	512.50	NIST Webbook
hvapt	131.70	kJ/mol	345.00	NIST Webbook
hvapt	135.00 ± 3.00	kJ/mol	431.50	NIST Webbook
psub	4.39e-08	kPa	329.02	Vapor and Sublimation Pressures of Three Normal Alkanes: C20, C24, and C28
psub	2.37e-08	kPa	326.15	Vapor and Sublimation Pressures of Three Normal Alkanes: C20, C24, and C28
psub	1.17e-08	kPa	323.05	Vapor and Sublimation Pressures of Three Normal Alkanes: C20, C24, and C28

pvap	1.70	kPa	550.62	Vapor-Liquid Equilibria of Binary Systems with Long-Chain Organic Compounds (Fatty Alcohol, Fatty Ester, Acylglycerol, and n-Paraffin) at Subatmospheric Pressures
pvap	0.04	kPa	473.64	A Gas Saturation Apparatus for Very Low Vapor or Sublimation Pressure Measurements (10 ⁻³ Pa): Vapor-Liquid Equilibria of n-Alkanes (n-C10, n-C24, n-C28)
pvap	0.01	kPa	453.28	A Gas Saturation Apparatus for Very Low Vapor or Sublimation Pressure Measurements (10 ⁻³ Pa): Vapor-Liquid Equilibria of n-Alkanes (n-C10, n-C24, n-C28)
pvap	0.16	kPa	494.80	A Gas Saturation Apparatus for Very Low Vapor or Sublimation Pressure Measurements (10 ⁻³ Pa): Vapor-Liquid Equilibria of n-Alkanes (n-C10, n-C24, n-C28)
pvap	2.93e-03	kPa	433.32	A Gas Saturation Apparatus for Very Low Vapor or Sublimation Pressure Measurements (10 ⁻³ Pa): Vapor-Liquid Equilibria of n-Alkanes (n-C10, n-C24, n-C28)

pvap	6.30e-04	kPa	413.50	A Gas Saturation Apparatus for Very Low Vapor or Sublimation Pressure Measurements (10 ⁻³ Pa): Vapor-Liquid Equilibria of n-Alkanes (n-C10, n-C24, n-C28)
pvap	6.51e-04	kPa	413.50	A Gas Saturation Apparatus for Very Low Vapor or Sublimation Pressure Measurements (10 ⁻³ Pa): Vapor-Liquid Equilibria of n-Alkanes (n-C10, n-C24, n-C28)
pvap	6.33e-04	kPa	413.50	A Gas Saturation Apparatus for Very Low Vapor or Sublimation Pressure Measurements (10 ⁻³ Pa): Vapor-Liquid Equilibria of n-Alkanes (n-C10, n-C24, n-C28)
pvap	1.26e-04	kPa	393.65	A Gas Saturation Apparatus for Very Low Vapor or Sublimation Pressure Measurements (10 ⁻³ Pa): Vapor-Liquid Equilibria of n-Alkanes (n-C10, n-C24, n-C28)
pvap	1.22e-04	kPa	393.65	A Gas Saturation Apparatus for Very Low Vapor or Sublimation Pressure Measurements (10 ⁻³ Pa): Vapor-Liquid Equilibria of n-Alkanes (n-C10, n-C24, n-C28)

pvap	1.80e-05	kPa	373.92	A Gas Saturation Apparatus for Very Low Vapor or Sublimation Pressure Measurements (10 ⁻³ Pa): Vapor-Liquid Equilibria of n-Alkanes (n-C10, n-C24, n-C28)
pvap	1.77e-05	kPa	373.92	A Gas Saturation Apparatus for Very Low Vapor or Sublimation Pressure Measurements (10 ⁻³ Pa): Vapor-Liquid Equilibria of n-Alkanes (n-C10, n-C24, n-C28)
pvap	2.14e-06	kPa	354.01	A Gas Saturation Apparatus for Very Low Vapor or Sublimation Pressure Measurements (10 ⁻³ Pa): Vapor-Liquid Equilibria of n-Alkanes (n-C10, n-C24, n-C28)
pvap	0.43	kPa	516.89	A Gas Saturation Apparatus for Very Low Vapor or Sublimation Pressure Measurements (10 ⁻³ Pa): Vapor-Liquid Equilibria of n-Alkanes (n-C10, n-C24, n-C28)
sfust	106.98	J/molxK	331.30	NIST Webbook
sfust	193.28	J/molxK	334.50	NIST Webbook

Pressure Dependent Properties

Property code	Value	Unit	Pressure [kPa]	Source
tbrp	551.20	K	2.00	NIST Webbook

Correlations

Information	Value
Property code	pvap
Equation	$\ln(P_{vp}) = A + B/(T + C)$
Coeff. A	1.69575e+01
Coeff. B	-7.53939e+03
Coeff. C	-9.37360e+01
Temperature range (K), min.	546.01
Temperature range (K), max.	741.12

Information	Value
Property code	pvap
Equation	$\ln(P_{vp}) = A + B/T + C \cdot \ln(T) + D \cdot T^2$
Coeff. A	3.55627e+02
Coeff. B	-3.27960e+04
Coeff. C	-4.75245e+01
Coeff. D	1.44719e-05
Temperature range (K), min.	481.15
Temperature range (K), max.	704.15

Sources

The Yaws Handbook of Vapor Pressure: <https://www.sciencedirect.com/book/9780128029992/the-yaws-handbook-of-vapor-pressure>

A Gas Saturation Apparatus for Very Low Vapor or Sublimation Pressure Measurements: <https://www.doi.org/10.1021/je7001122>

Fluorinated Alkanes: Effects of Alkane Chain Length and Equivalents of Binary Carbon Surfaces with Long-Chain Organic Compounds: <https://www.doi.org/10.1016/j.fluid.2004.10.021>

Measurement of Enthalpy Curves of Phase Change Materials via DSC and Crippen Method: <https://www.doi.org/10.1021/acs.jced.8b00168>

Measurement of Enthalpy Curves of Phase Change Materials via DSC and Crippen Method: <https://www.doi.org/10.1021/acs.jced.7b00363>

Measurement of Enthalpy Curves of Phase Change Materials via DSC and Crippen Method: <https://www.doi.org/10.1016/j.fluid.2014.07.017>

Measurement of Enthalpy Curves of Phase Change Materials via DSC and Crippen Method: <https://www.cheric.org/files/research/kdb/mol/mol28.mol>

Joback Method: https://en.wikipedia.org/wiki/Joback_method

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

KDB Vapor Pressure Data: <https://www.cheric.org/research/kdb/hcprop/showprop.php?cmpid=28>

Measurement of enthalpy curves of phase change materials via DSC and Crippen Method: <https://www.doi.org/10.1016/j.tca.2014.09.022>

Measurement of enthalpy curves of phase change materials via DSC and Crippen Method: https://www.chemeo.com/doc/models/crippen_log10ws

Measurement of enthalpy curves of phase change materials via DSC and Crippen Method: <https://www.doi.org/10.1021/acs.jced.8b01139>

Measurement of enthalpy curves of phase change materials via DSC and Crippen Method: <https://www.doi.org/10.1021/acs.jced.8b01139>

Vapor Pressures and Vaporization Enthalpies of the n-Alkanes from C₂₁ to C₃₀ at 298.15 K by Correlation Gas Chromatography: McGowan Method.

<https://www.doi.org/10.1021/je0301747>

Vapor and Sublimation Pressures of Three Normal Alkanes: C₂₀, C₂₄, and C₂₈. Thermal conductivity of heavy, even-carbon number n-alkanes (C₂₂ to C₃₂). Comparison of Results by Correlation Gas Chromatography with Another Gas Chromatographic Retention Time Technique. The Effects of Retention Time Coincidence on Vaporization Enthalpy and Vapor Pressure.

<http://webbook.nist.gov/cgi/cbook.cgi?ID=C630024&Units=SI>

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

<https://www.doi.org/10.1021/je800534x>

<https://www.doi.org/10.1016/j.fluid.2018.08.016>

<https://www.doi.org/10.1021/acs.jced.5b00444>

Legend

cp_g:	Ideal gas heat capacity
cp_l:	Liquid phase heat capacity
dv_{isc}:	Dynamic viscosity
gf:	Standard Gibbs free energy of formation
hf:	Enthalpy of formation at standard conditions
hf_{us}:	Enthalpy of fusion at standard conditions
hf_{ust}:	Enthalpy of fusion at a given temperature
h_{sub}:	Enthalpy of sublimation at standard conditions
h_{subt}:	Enthalpy of sublimation at a given temperature
h_{vap}:	Enthalpy of vaporization at standard conditions
h_{vapt}:	Enthalpy of vaporization at a given temperature
log₁₀w_s:	Log ₁₀ of Water solubility in mol/l
log_p:	Octanol/Water partition coefficient
mc_{vol}:	McGowan's characteristic volume
pc:	Critical Pressure
ps_{ub}:	Sublimation pressure
pv_{ap}:	Vapor pressure
rin_{pol}:	Non-polar retention indices
sf_{ust}:	Entropy of fusion at a given temperature
tb:	Normal Boiling Point Temperature
tb_{rp}:	Boiling point at reduced pressure
tc:	Critical Temperature
tf:	Normal melting (fusion) point
tt:	Triple Point Temperature
vc:	Critical Volume

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