

# 14-Tetracosene, 27-methyl

**InChI:** InChI=1S/C41H82/c1-4-6-8-10-12-14-16-17-18-19-20-21-22-23-24-25-26-27-28-30-32-34-36-38-40-41(3)39-37-35-33-31-29-15-13-11-9-7-5-2/h22-23,41H,4-21,24-40H2,1-3H3/b23-22+

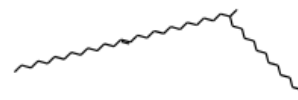
**InChI Key:** KDWZUEQMMIVDOK-GHVJWSGMSA-N

**Formula:** C41H82

**SMILES:**

CCCCCCCCCCCCC=CCCCCCCCCCCCC(C)CCCCCCCCCCCCC

**Molecular Weight:** 575.09



## Physical Properties

Property	Value	Unit	Source
$\Delta_f G^\circ$	372.12	kJ/mol	Joback Method
$\Delta_f H^\circ_{\text{gas}}$	-777.63	kJ/mol	Joback Method
$\Delta_{\text{fus}} H^\circ$	98.62	kJ/mol	Joback Method
$\Delta_{\text{vap}} H^\circ$	106.43	kJ/mol	Joback Method
$\log P_{\text{oct/wat}}$	15.87		Crippen Method
$P_c$	385.37	kPa	Joback Method
$T_{\text{boil}}$	1141.20	K	Joback Method
$T_c$	1526.95	K	Joback Method
$T_{\text{fus}}$	531.75	K	Joback Method
$V_c$	2.31	m <sup>3</sup> /kg-mol	Joback Method

## Temperature Dependent Properties

Property	Value	Unit	Temperature (K)	Source
$C_{p,\text{gas}}$	2219.84	J/mol×K	1141.2	Joback Method
$\eta$	0.00	Paxs	1141.2	Joback Method

## Sources

**Joback Method:** [https://en.wikipedia.org/wiki/Joback\\_method](https://en.wikipedia.org/wiki/Joback_method)

**NIST Webbook:** [http://webbook.nist.gov/cgi/inchi/InChI=1S/C41H82/c1-4-6-8-10-12-14-16-17-18-19-20-21-22-23-24-25-26-27-28-30-32-34-36-38-40-41\(3\)39-37-35-33-31-29-15-13-11-9-7-5-2/h22-23,41H,4-21,24-40H2,1-3H3/b23-22+](http://webbook.nist.gov/cgi/inchi/InChI=1S/C41H82/c1-4-6-8-10-12-14-16-17-18-19-20-21-22-23-24-25-26-27-28-30-32-34-36-38-40-41(3)39-37-35-33-31-29-15-13-11-9-7-5-2/h22-23,41H,4-21,24-40H2,1-3H3/b23-22+)

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

## Legend

$C_{p,gas}$ : Ideal gas heat capacity (J/molxK).

$\eta$ : Dynamic viscosity (Pa $\times$ s).

$\Delta_f G^\circ$ : Standard Gibbs free energy of formation (kJ/mol).

$\Delta_f H^\circ_{gas}$ : Enthalpy of formation at standard conditions (kJ/mol).

$\Delta_{fus} H^\circ$ : Enthalpy of fusion at standard conditions (kJ/mol).

$\Delta_{vap} H^\circ$ : Enthalpy of vaporization at standard conditions (kJ/mol).

$logP_{oct/wat}$ : Octanol/Water partition coefficient .

$P_c$ : Critical Pressure (kPa).

$T_{boil}$ : Normal Boiling Point Temperature (K).

$T_c$ : Critical Temperature (K).

$T_{fus}$ : Normal melting (fusion) point (K).

$V_c$ : Critical Volume (m<sup>3</sup>/kg-mol).

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