

# 5-Tetradecene, (E)-

<b>Other names:</b>	(5E)-5-Tetradecene (E)-5-Tetradecene E-Tetradec-5-ene trans-5-Tetradecene
<b>Inchi:</b>	InChI=1S/C14H28/c1-3-5-7-9-11-13-14-12-10-8-6-4-2/h9,11H,3-8,10,12-14H2,1-2H3/b1
<b>InchiKey:</b>	SNIFAVVHRQZYGO-PKNBQFBNSA-N
<b>Formula:</b>	C14H28
<b>SMILES:</b>	CCCC=CCCCCCCC
<b>Mol. weight [g/mol]:</b>	196.37
<b>CAS:</b>	41446-66-6

## Physical Properties

Property code	Value	Unit	Source
gf	147.22	kJ/mol	Joback Method
hf	-215.07	kJ/mol	Joback Method
hfus	32.22	kJ/mol	Joback Method
hvap	46.72	kJ/mol	Joback Method
log10ws	-5.54		Crippen Method
logp	5.483		Crippen Method
mcvol	203.820	ml/mol	McGowan Method
pc	1587.28	kPa	Joback Method
rinpol	1380.00		NIST Webbook
rinpol	1378.00		NIST Webbook
rinpol	1391.00		NIST Webbook
rinpol	1386.00		NIST Webbook
rinpol	1387.00		NIST Webbook
rinpol	1391.00		NIST Webbook
rinpol	1378.00		NIST Webbook
rinpol	1377.80		NIST Webbook
rinpol	1378.00		NIST Webbook
rinpol	1378.00		NIST Webbook
ripol	1437.00		NIST Webbook
ripol	1427.10		NIST Webbook
ripol	1427.80		NIST Webbook
ripol	1419.00		NIST Webbook
ripol	1425.00		NIST Webbook
ripol	1426.00		NIST Webbook

ripol	1426.00		NIST Webbook
ripol	1421.00		NIST Webbook
ripol	1426.00		NIST Webbook
ripol	1426.00		NIST Webbook
ripol	1427.00		NIST Webbook
ripol	1421.00		NIST Webbook
ripol	1428.00		NIST Webbook
ripol	1435.00		NIST Webbook
ripol	1436.00		NIST Webbook
ripol	1436.00		NIST Webbook
ripol	1421.50		NIST Webbook
ripol	1438.00		NIST Webbook
ripol	1430.00		NIST Webbook
ripol	1431.00		NIST Webbook
ripol	1432.00		NIST Webbook
ripol	1432.00		NIST Webbook
ripol	1433.00		NIST Webbook
ripol	1418.00		NIST Webbook
ripol	1425.60		NIST Webbook
ripol	1419.80		NIST Webbook
ripol	1426.10		NIST Webbook
ripol	1415.80		NIST Webbook
ripol	1419.30		NIST Webbook
ripol	1425.00		NIST Webbook
ripol	1425.70		NIST Webbook
ripol	1418.60		NIST Webbook
ripol	1420.60		NIST Webbook
ripol	1425.60		NIST Webbook
ripol	1426.90		NIST Webbook
ripol	1435.00		NIST Webbook
ripol	1436.00		NIST Webbook
ripol	1437.00		NIST Webbook
ripol	1438.00		NIST Webbook
ripol	1421.50		NIST Webbook
ripol	1426.00		NIST Webbook
ripol	1427.00		NIST Webbook
ripol	1435.00		NIST Webbook
ripol	1432.00		NIST Webbook
ripol	1419.80		NIST Webbook
ripol	1425.70		NIST Webbook
ripol	1436.00		NIST Webbook
tb	523.88	K	Joback Method
tc	689.82	K	Joback Method
tf	242.46	K	Joback Method

## Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	488.90	J/molxK	523.88	Joback Method
cpg	571.08	J/molxK	662.16	Joback Method
cpg	556.03	J/molxK	634.51	Joback Method
cpg	540.31	J/molxK	606.85	Joback Method
cpg	523.90	J/molxK	579.19	Joback Method
cpg	506.77	J/molxK	551.54	Joback Method
cpg	585.50	J/molxK	689.82	Joback Method
dvisc	0.0001553	Paxs	523.88	Joback Method
dvisc	0.0002098	Paxs	476.98	Joback Method
dvisc	0.0003029	Paxs	430.07	Joback Method
dvisc	0.0004782	Paxs	383.17	Joback Method
dvisc	0.0008577	Paxs	336.27	Joback Method
dvisc	0.0018593	Paxs	289.36	Joback Method
dvisc	0.0054364	Paxs	242.46	Joback Method

## Sources

<b>NIST Webbook:</b>	<a href="http://webbook.nist.gov/cgi/cbook.cgi?ID=C41446666&amp;Units=SI">http://webbook.nist.gov/cgi/cbook.cgi?ID=C41446666&amp;Units=SI</a>
<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.chemeo.com/doc/models/crippen_log10ws">https://www.chemeo.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>

## 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>h<sub>vap</sub>:</b>	Enthalpy of vaporization at standard conditions
<b>log<sub>10</sub>w<sub>s</sub>:</b>	Log <sub>10</sub> of Water solubility in mol/l
<b>log<sub>p</sub>:</b>	Octanol/Water partition coefficient
<b>m<sub>cvol</sub>:</b>	McGowan's characteristic volume
<b>p<sub>c</sub>:</b>	Critical Pressure
<b>r<sub>inpol</sub>:</b>	Non-polar retention indices
<b>r<sub>ipol</sub>:</b>	Polar retention indices
<b>t<sub>b</sub>:</b>	Normal Boiling Point Temperature
<b>t<sub>c</sub>:</b>	Critical Temperature
<b>t<sub>f</sub>:</b>	Normal melting (fusion) point
<b>v<sub>c</sub>:</b>	Critical Volume

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