

# I-Leucine, N-allyloxycarbonyl-N-methyl-, pentadecyl ester

<b>Inchi:</b>	InChI=1S/C26H49NO4/c1-6-8-9-10-11-12-13-14-15-16-17-18-19-21-30-25(28)24(22-23)
<b>InchiKey:</b>	XOEJOXISVWTHCE-UHFFFAOYSA-N
<b>Formula:</b>	C26H49NO4
<b>SMILES:</b>	C=CCOC(=O)N(C)C(CC(C)C)C(=O)OCCCCCCCCCCCCCCC
<b>Mol. weight [g/mol]:</b>	439.67

## Physical Properties

Property code	Value	Unit	Source
gf	-106.06	kJ/mol	Joback Method
hf	-887.17	kJ/mol	Joback Method
hfus	63.37	kJ/mol	Joback Method
hvap	92.38	kJ/mol	Joback Method
log10ws	-7.70		Crippen Method
logp	7.290		Crippen Method
mcvol	397.760	ml/mol	McGowan Method
pc	786.39	kPa	Joback Method
rinpol	2807.00		NIST Webbook
rinpol	2807.00		NIST Webbook
tb	955.10	K	Joback Method
tc	1174.45	K	Joback Method
tf	527.81	K	Joback Method
vc	1.526	m <sup>3</sup> /kmol	Joback Method

## Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	1361.12	J/mol×K	955.10	Joback Method
cpg	1381.83	J/mol×K	991.66	Joback Method
cpg	1400.92	J/mol×K	1028.22	Joback Method
cpg	1418.46	J/mol×K	1064.78	Joback Method
cpg	1434.50	J/mol×K	1101.33	Joback Method
cpg	1449.13	J/mol×K	1137.89	Joback Method
cpg	1462.40	J/mol×K	1174.45	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=U321906&amp;Units=SI">http://webbook.nist.gov/cgi/cbook.cgi?ID=U321906&amp;Units=SI</a>

# Legend

<b>cpg:</b>	Ideal gas heat capacity
<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>hvp:</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>rinp:</b>	Non-polar retention indices
<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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