

# Pheneturide

<b>Other names:</b>	Benzeneacetamide, N-(aminocarbonyl)-«alpha»-ethyl- Urea, (2-phenylbutyryl)- Benuride Ethylphenacemide 1-((Ethyl)phenylacetyl)urea Lircapyl M 551 Phenuride N-(«alpha»-Phenylbutyryl)urea 2-Phenylbutyrylurea Phenylethylacetylurea Phenylethylacetyluree S 46 EPA PBU
<b>Inchi:</b>	InChI=1S/C11H14N2O2/c1-2-9(10(14)13-11(12)15)8-6-4-3-5-7-8/h3-7,9H,2H2,1H3,(H3,1
<b>InchiKey:</b>	AJOQSQHYDOFIOX-UHFFFAOYSA-N
<b>Formula:</b>	C11H14N2O2
<b>SMILES:</b>	CCC(C(=O)NC(N)=O)c1ccccc1
<b>Mol. weight [g/mol]:</b>	206.24
<b>CAS:</b>	90-49-3

## Physical Properties

Property code	Value	Unit	Source
gf	49.71	kJ/mol	Joback Method
hf	-177.02	kJ/mol	Joback Method
hfus	28.26	kJ/mol	Joback Method
hvap	72.54	kJ/mol	Joback Method
log10ws	-2.65		Crippen Method
logp	1.375		Crippen Method
mcvol	165.190	ml/mol	McGowan Method
pc	3306.75	kPa	Joback Method
rinpol	1465.00		NIST Webbook
rinpol	1465.00		NIST Webbook
tb	707.76	K	Joback Method
tc	937.06	K	Joback Method
tf	460.93	K	Joback Method

vc

0.614

m<sup>3</sup>/kmol

Joback Method

## Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	446.15	J/mol×K	707.76	Joback Method
cpg	458.69	J/mol×K	745.98	Joback Method
cpg	470.26	J/mol×K	784.19	Joback Method
cpg	480.90	J/mol×K	822.41	Joback Method
cpg	490.68	J/mol×K	860.62	Joback Method
cpg	499.63	J/mol×K	898.84	Joback Method
cpg	507.82	J/mol×K	937.06	Joback Method

## Sources

**Joback Method:**

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

**McGowan Method:**

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

**NIST Webbook:**

<http://webbook.nist.gov/cgi/cbook.cgi?ID=C90493&Units=SI>

**Crippen Method:**

<http://pubs.acs.org/doi/abs/10.1021/ci990307l>

**Crippen Method:**

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

## 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>hvap:</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>rinpol:</b>	Non-polar retention indices
<b>tb:</b>	Normal Boiling Point Temperature
<b>tc:</b>	Critical Temperature

**tf:** Normal melting (fusion) point

**vc:** Critical Volume

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