

# Fenproporex-M (desamino-oxo-HO-methoxy), AC

Inchi:	InChI=1S/C12H14O4/c1-8(13)6-10-4-5-11(16-9(2)14)7-12(10)15-3/h4-5,7H,6H2,1-3H3
InchiKey:	SOLHVFDPMCUTMO-UHFFFAOYSA-N
Formula:	C12H14O4
SMILES:	COc1cc(OC(C)=O)ccc1CC(C)=O
Mol. weight [g/mol]:	222.24

## Physical Properties

Property code	Value	Unit	Source
gf	-324.53	kJ/mol	Joback Method
hf	-567.02	kJ/mol	Joback Method
hfus	25.67	kJ/mol	Joback Method
hvap	64.22	kJ/mol	Joback Method
log10ws	-2.41		Crippen Method
logp	1.752		Crippen Method
mcvol	171.060	ml/mol	McGowan Method
pc	2558.51	kPa	Joback Method
rinqol	1600.00		NIST Webbook
tb	663.18	K	Joback Method
tc	875.86	K	Joback Method
tf	420.78	K	Joback Method
vc	0.647	m <sup>3</sup> /kmol	Joback Method

## Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	435.52	J/molxK	663.18	Joback Method
cpg	448.70	J/molxK	698.63	Joback Method
cpg	461.11	J/molxK	734.07	Joback Method
cpg	472.74	J/molxK	769.52	Joback Method
cpg	483.58	J/molxK	804.97	Joback Method
cpg	493.62	J/molxK	840.42	Joback Method
cpg	502.86	J/molxK	875.86	Joback Method
dvisc	0.0009239	Paxs	420.78	Joback Method
dvisc	0.0005957	Paxs	461.18	Joback Method

dvisc	0.0004122	Paxs	501.58	Joback Method
dvisc	0.0003014	Paxs	541.98	Joback Method
dvisc	0.0002301	Paxs	582.38	Joback Method
dvisc	0.0001820	Paxs	622.78	Joback Method
dvisc	0.0001481	Paxs	663.18	Joback Method

## Sources

<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=R275106&amp;Units=SI">http://webbook.nist.gov/cgi/cbook.cgi?ID=R275106&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>

## 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>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
<b>tf:</b>	Normal melting (fusion) point
<b>vc:</b>	Critical Volume

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