

o-anisic acid, pent-2-en-4-ynyl ester

Inchi:

InchiKey:

Formula:

SMILES:

Mol. weight [g/mol]:

InChI=1S/C13H12O3/c1-3-4-7-10-16-13(14)11-8-5-6-9-12(11)15-2/h1,4-9H,10H2,2H3/b7

MRQQDBRYOKWHPC-QPJXVBHSA-N

C13H12O3

C#C/C=C/COC(=O)c1ccccc1OC

216.24

Physical Properties

Property code	Value	Unit	Source
hf	-132.64	kJ/mol	Cheméo Relay (1.0)
hfus	30.23	kJ/mol	Joback Method
hvap	83.29	kJ/mol	Cheméo Relay (1.0)
log10ws	-3.75		Cheméo Relay (1.0)
logp	2.041		Crippen Method
mcvol	170.680	ml/mol	McGowan Method
pc	2709.85	kPa	Joback Method
rinpol	1751.20		NIST Webbook
rinpol	1751.20		NIST Webbook
tb	572.51	K	Cheméo Relay (1.0)
tf	333.36	K	Cheméo Relay (1.0)

Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	405.39	J/mol×K	621.49	Joback Method
cpg	418.89	J/mol×K	658.83	Joback Method
cpg	431.53	J/mol×K	696.18	Joback Method
cpg	443.34	J/mol×K	733.52	Joback Method
cpg	454.36	J/mol×K	770.87	Joback Method
cpg	464.61	J/mol×K	808.21	Joback Method
cpg	474.12	J/mol×K	845.55	Joback Method

Sources

McGowan Method:	http://link.springer.com/article/10.1007/BF02311772
Crippen Method:	http://pubs.acs.org/doi/abs/10.1021/ci990307l
Crippen Method:	https://www.chemeo.com/doc/models/crippen_log10ws
Cheméo Relay (1.0):	
Cheméo Relay (1.0, beta):	
NIST Webbook:	http://webbook.nist.gov/cgi/cbook.cgi?ID=U292573&Units=SI
Joback Method:	https://en.wikipedia.org/wiki/Joback_method

Legend

cpg:	Ideal gas heat capacity
hf:	Enthalpy of formation at standard conditions
hfus:	Enthalpy of fusion at standard conditions
hvap:	Enthalpy of vaporization at standard conditions
log10ws:	Log10 of Water solubility in mol/l
logp:	Octanol/Water partition coefficient
mcvol:	McGowan's characteristic volume
pc:	Critical Pressure
rinpol:	Non-polar retention indices
tb:	Normal Boiling Point Temperature
tf:	Normal melting (fusion) point

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