

# CPIB PFB ester

<b>Other names:</b>	Clofibric acid, PFB
<b>Inchi:</b>	InChI=1S/C17H12ClF5O3/c1-17(2,26-9-5-3-8(18)4-6-9)16(24)25-7-10-11(19)13(21)15(23)
<b>InchiKey:</b>	LWCTUOXYEKQVKP-UHFFFAOYSA-N
<b>Formula:</b>	C17H12ClF5O3
<b>SMILES:</b>	CC(C)(Oc1ccc(Cl)cc1)C(=O)OCc1c(F)c(F)c(F)c(F)c1F
<b>Mol. weight [g/mol]:</b>	394.72

## Physical Properties

Property code	Value	Unit	Source
gf	-1062.76	kJ/mol	Joback Method
hf	-1372.03	kJ/mol	Joback Method
hfus	41.69	kJ/mol	Joback Method
hvap	72.53	kJ/mol	Joback Method
log10ws	-6.69		Crippen Method
logp	4.936		Crippen Method
mcvol	237.270	ml/mol	McGowan Method
pc	1610.29	kPa	Joback Method
rinpol	1934.00		NIST Webbook
rinpol	1933.00		NIST Webbook
ripol	2525.00		NIST Webbook
ripol	2525.00		NIST Webbook
ripol	2603.00		NIST Webbook
tb	800.86	K	Joback Method
tc	1005.98	K	Joback Method
tf	538.99	K	Joback Method
vc	0.942	m3/kmol	Joback Method

## Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	657.84	J/molxK	800.86	Joback Method
cpg	669.29	J/molxK	835.05	Joback Method
cpg	679.83	J/molxK	869.23	Joback Method
cpg	689.50	J/molxK	903.42	Joback Method

cpg	698.30	J/mol×K	937.61	Joback Method
cpg	706.25	J/mol×K	971.79	Joback Method
cpg	713.38	J/mol×K	1005.98	Joback Method

## Sources

<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=R14032&amp;Units=SI">http://webbook.nist.gov/cgi/cbook.cgi?ID=R14032&amp;Units=SI</a>
<b>Crippen Method:</b>	<a href="http://pubs.acs.org/doi/abs/10.1021/ci990307l">http://pubs.acs.org/doi/abs/10.1021/ci990307l</a>
<b>Crippen Method:</b>	<a href="https://www.chemeo.com/doc/models/crippen_log10ws">https://www.chemeo.com/doc/models/crippen_log10ws</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>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>ripol:</b>	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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