

PHOSPHINE, TRIS(4-METHOXYPHENYL)-

Other names:	NSC 136458
	Phosphine, tris(p-methoxyphenyl)-
	Phosphorus tri-(p-methoxyphenyl)
	Tri(p-methoxyphenyl)phosphine
	Tri-p-anisylphosphine
	Trianisylphosphine
	Tris(4-anisyl)phosphine
	Tris(p-anisyl)phosphine
	tri-(4-Methoxyphenyl)phosphine
	tris(4-methoxyphenyl)phosphane
	tris(4-methoxyphenyl)phosphine
	tris(p-methoxyphenyl)phosphine
Inchi:	InChI=1S/C21H21O3P/c1-22-16-4-10-19(11-5-16)25(20-12-6-17(23-2)7-13-20)21-14-8-1
InchiKey:	UYUUAUOYLFIRJG-UHFFFAOYSA-N
Formula:	C21H21O3P
SMILES:	COc1ccc(P(c2ccc(OC)cc2)c2ccc(OC)cc2)cc1
Mol. weight [g/mol]:	352.37

Physical Properties

Property code	Value	Unit	Source
hf	-249.60	kJ/mol	Cheméo Relay (1.0)
hfus	25.67	kJ/mol	Solubilities of Phosphorus-Containing Compounds in Selected Solvents
hvap	110.12	kJ/mol	Cheméo Relay (1.0)
ie	7.48	eV	NIST Webbook
log10ws	-6.92		Cheméo Relay (1.0)
logp	3.471		Crippen Method
mcvol	273.540	ml/mol	McGowan Method
pc	1932.80	kPa	Cheméo Relay (1.0, beta)
tb	727.02	K	Cheméo Relay (1.0)
tc	972.78	K	Cheméo Relay (1.0)

tf	403.65	K	Solubilities of Phenylphosphinic Acid, Hydroxymethylphenylphosphinic Acid, p-Methoxyphenylphosphinic Acid, p-Methoxyphenylhydroxymethylphosphinic Acid, Triphenylphosphine, Tri(p-methoxyphenyl)phosphine, and Tri(p-methoxyphenyl)phosphine Oxide in Selected Solvents
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Sources

Solubilities of Phenylphosphinic Acid, Hydroxymethylphenylphosphinic Acid, p-Methoxyphenylphosphinic Acid, p-Methoxyphenylhydroxymethylphosphinic Acid, Triphenylphosphine, Tri(p-methoxyphenyl)phosphine, and Tri(p-methoxyphenyl)phosphine Oxide in Selected Solvents: <https://www.doi.org/10.1021/je050377q>
 Solubilities of Phenylphosphinic Acid, Hydroxymethylphenylphosphinic Acid, p-Methoxyphenylphosphinic Acid, p-Methoxyphenylhydroxymethylphosphinic Acid, Triphenylphosphine, Tri(p-methoxyphenyl)phosphine, and Tri(p-methoxyphenyl)phosphine Oxide in Selected Solvents: <https://www.doi.org/10.1021/je100341q>
 NIST Webbook: <http://webbook.nist.gov/cgi/cbook.cgi?ID=C855389&Units=SI>
 Crippen Method: <http://link.springer.com/article/10.1007/BF02311772>
 Cheméo Relay (1.0): <http://pubs.acs.org/doi/abs/10.1021/ci990307l>
 Cheméo Relay (1.0, beta): https://www.chemeo.com/doc/models/crippen_log10ws

Legend

hf: Enthalpy of formation at standard conditions
hfus: Enthalpy of fusion at standard conditions
hvac: Enthalpy of vaporization at standard conditions
ie: Ionization energy
log10ws: Log10 of Water solubility in mol/l
logp: Octanol/Water partition coefficient
mcvol: McGowan's characteristic volume
pc: Critical Pressure
tb: Normal Boiling Point Temperature
tc: Critical Temperature
tf: Normal melting (fusion) point

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