User: Fakhouri, Morgan
Westchester Dept. of Labs & Research
Date: 12/11/2001 Time: 11:25:44

Sample: AD27547
Analysis: $B1525 Organic Chemicals - 525.2; lab blank
Units: ug
Started: 11/16/01 10:58 Ended: 12/5/01 10:58 Analyst: MF
Result source: manual entry
Page: 1

	Component Name	Viol	Result	Qualifier	MDL
1	Hexachlorocyclopentadiene		< MDL		0.10
2	Hexachlorobenzene		< MDL		0.10
3	Simazine		< MDL		0.07
4	Atrazine		< MDL		0.20
5	Alachlor		< MDL		0.20
6	bis(2-Ethylhexyl)adipate		< MDL		0.60
7	bis(2-Ethylhexyl)phthalate		< MDL		0.60
8	Benzo(a)pyrene		< MDL		0.20
9	EPTC		< MDL		1.0
10	2,6-Dinitrotoluene		< MDL		2.0
11	2,4-Dinitrotoluene		< MDL		2.0
12	Molinate		< MDL		0.90
13	Terbacil		< MDL		2.0
14	Acetochlor		< MDL		2.0
15	Metribuzin		< MDL		0.15
16	Metolachlor		< MDL		0.75
17	Butachlor		< MDL		0.40
18	4,4-DDE		< MDL		0.80
19	[redacted]				
20	Surr_Triphenyl phosphate		5.30		
21	Surr_Perylene-d12		4.19		

3.5 - 6.5

Quantitation Report (QT Reviewed)

Data File : C:\MSDCHEM\1\DATA\120501\1116LB.D Vial: 5
Acq On : 5 Dec 2001 6:24 pm Operator: McLean
ample : 11/16 run Inst : Instrumen
Misc : Multiplr: 1.00
MS Integration Params: BENZO.P
Quant Time: Dec 11 09:10:34 2001 Quant Results File: 120501.RES

Quant Method : C:\MSDCHEM\1\5973N\120501.M (RTE Integrator)
Title : Quantitation For Method 525.2
Last Update : Thu Dec 06 10:31:04 2001
Response via : Initial Calibration
DataAcq Meth : 120501

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Acenaphthene-d10	18.31	164	4810429	5.00	ug/L	0.00
9) Phenanthrene D-10	22.64	188	7400763	5.00	ug/L	0.00
20) Chrysene D-12	30.42	240	5218776	5.00	ug/L	0.00
System Monitoring Compounds						
2) 1,3-Dimethyl-2-nitrobenzen	13.23	134	936795	3.67	ug/L	0.00
Spiked Amount 5.000			Recovery = 73.40%			
19) Triphenyl Phosphate surr	29.69	326	1359468	5.30	ug/L	0.00
Spiked Amount 5.000			Recovery = 106.00%			
21) Perylene D-12 surr	34.44	264	3066374	4.19	ug/L	0.00
Spiked Amount 5.000			Recovery = 83.80%			
Target Compounds						Qvalue
3) Hexachlorocyclopentadiene	0.00	237	0	N.D.		
4) Hexachlorobenzene	0.00	284	0	N.D.		
5) EPTC	0.00	128	0	N.D.		
6) 2,6-Dinitrotoluene	0.00	165	0	N.D.	d	
7) 2,4-Dinitotoluene	0.00	165	0	N.D.	d	
8) Molinate	19.09	126	9997	N.D.		
10) Simazine	0.00	201	0	N.D.		
11) Atrazine	0.00	200	0	N.D.		
12) Alachlor	0.00	160	0	N.D.		
13) Terbacil	0.00	161	0	N.D.	d	
14) Acetochlor	0.00	146	0	N.D.		
15) Metribuzin	0.00	198	0	N.D.		
16) Metolachlor	0.00	162	0	N.D.		
17) Butachlor	0.00	160	0	N.D.	d	
18) 4,4'-DDE	0.00	246	0	N.D.	d	
22) Bis (2-Ethylhexyl) adipate	29.62	129	9060	0.12	ug/L #	11
23) Bis (2-Ethylhexyl) phthala	31.07	149	120215	0.26	ug/L	100
24) Benzo (a) pyrene	0.00	252	0	N.D.	d	

(#) = qualifier out of range (m) = manual integration
1116LB.D 120501.M Tue Dec 11 09:13:21 2001 Page 1

Data File : C:\MSDCHEM\1\DATA\120501\1116LB.D Vial: 5
Acq On : 5 Dec 2001 6:24 pm Operator: McLean
Sample : 11/16 run Inst : Instrumen
Misc : Multiplr: 1.00
MS Integration Params: BENZO.P
Quant Time: Dec 11 9:13 2001 Quant Results File: 120501.RES

Method : C:\MSDCHEM\1\5973N\120501.M (RTE Integrator)
Title : Quantitation For Method 525.2
Last Update : Thu Dec 06 10:31:04 2001
Response via : Initial Calibration