User: Fakhouri, Morgan
Westchester Dept. of Labs & Research
Date: 12/20/2001 Time: 11:59:34

Sample: AD28847
Analysis: $B1525 Organic Chemicals - 525.2; lab blank
Units: ug
Started: 12/20/01 11:54 Ended: 12/20/01 11:54 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	Surr_1,3-Dimethyl-2-nitrobenzen		4.89		
20	Surr_Triphenyl phosphate		4.71		
21	Surr_Perylene-d12		5.02		

Quantitation Report (QT Reviewed)

Data File : C:\MSDCHEM\1\DATA\121701\1127LBA.D Vial: 27
Acq On : 17 Dec 2001 4:16 pm Operator: McLean
Sample : Inst : Instrumen
Misc : Multiplr: 1.00
MS Integration Params: BENZO.P
Quant Time: Dec 17 16:53:23 2001 Quant Results File: 121401.RES

Quant Method : C:\MSDCHEM\1\5973N\121401.M (RTE Integrator)
Title : Quantitation For Method 525.2
Last Update : Mon Dec 17 15:06:05 2001
Response via : Initial Calibration
DataAcq Meth : 121401

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Acenaphthene-d10	18.20	164	5385816	5.00	ug/L	0.00
9) Phenanthrene D-10	22.53	188	8319552	5.00	ug/L	0.00
20) Chrysene D-12	30.30	240	5299936	5.00	ug/L	0.00
System Monitoring Compounds						
2) 1,3-Dimethyl-2-nitrobenzen	13.13	134	1348111	4.89	ug/L	0.00
Spiked Amount 5.000			Recovery = 97.80%			
19) Triphenyl Phosphate surr	29.58	326	1192866	4.71	ug/L	0.00
Spiked Amount 5.000			Recovery = 94.20%			
21) Perylene D-12 surr	34.31	264	3520740	5.02	ug/L	0.00
Spiked Amount 5.000			Recovery = 100.40%			
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	18.94	165	1119	0.31	ug/L #	21
8) Molinate	0.00	126	0	N.D. 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	29.51	129	9041	0.39	ug/L #	31
23) Bis (2-Ethylhexyl) phthala	30.95	149	136786	0.46	ug/L	100
24) Benzo (a) pyrene	0.00	252	0	N.D. d		

(#) = qualifier out of range (m) = manual integration

Quantitation Report (QT Reviewed)

Data File : C:\MSDCHEM\1\DATA\121701\1127LBA.D Vial: 27
Acq On : 17 Dec 2001 4:16 pm Operator: McLean
Sample : Inst : Instrumen
Misc : Multiplr: 1.00
MS Integration Params: BENZO.P
Quant Time: Dec 20 11:58 2001 Quant Results File: 121401.RES

Method : C:\MSDCHEM\1\5973N\121401.M (RTE Integrator)
Title : Quantitation For Method 525.2
Last Update : Tue Dec 18 09:35:41 2001
Response via : Initial Calibration