User: Fakhouri, Morgan
Westchester Dept. of Labs & Research
Date: 12/12/2001 Time: 16:00:22

Sample: AD27232
Analysis: $B1525 Organic Chemicals - 525.2; lab blank
Units: ug
Started: 11/15/01 15:57 Ended: 12/6/01 15:57 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					
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					
20	Surr_Triphenyl phosphate		4.39		
21	Surr_Perylene-d12		3.06		

Quantitation Report (QT Reviewed)

Data File : C:\MSDCHEM\1\DATA\120601\1114LBK.D Vial: 29
Acq On : 7 Dec 2001 2:19 pm Operator: McLean
Sample : Inst : Instrumen
Misc : Multiplr: 1.00
MS Integration Params: BENZO.P
Quant Time: Dec 12 11:55:38 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	18.32	164	3920592	5.00	ug/L	0.00
9) Phenanthrene D-10	22.65	188	6784721	5.00	ug/L	0.00
20) Chrysene D-12	30.43	240	4002597	5.00	ug/L	0.00
System Monitoring Compounds						
2) 1,3-Dimethyl-2-nitrobenzen	13.23	134	779452	3.75	ug/L	0.00
Spiked Amount 5.000			Recovery = 75.00%			
19) Triphenyl Phosphate surr	29.70	326	1032641	4.39	ug/L	0.01
Spiked Amount 5.000			Recovery = 87.80%			
21) Perylene D-12 surr	34.46	264	1717757	3.06	ug/L	0.02
Spiked Amount 5.000			Recovery = 61.20%			
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	19.08	126	7014	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	26.77	160	28354	0.10	ug/L #	7
18) 4,4'-DDE	27.52	246	163950	0.58	ug/L #	23
22) Bis (2-Ethylhexyl) adipate	29.62	129	11870	0.13	ug/L #	20
23) Bis (2-Ethylhexyl) phthala	31.07	149	541968	0.70	ug/L	92
24) Benzo (a) pyrene	0.00	252	0	N.D.		

(#) = qualifier out of range (m) = manual integration
1114LBK.D 120501.M Wed Dec 12 11:56:22 2001 Page 1

Quantitation Report (QT Reviewed)

Data File : C:\MSDCHEM\1\DATA\120601\1114LBK.D Vial: 29
Acq On : 7 Dec 2001 2:19 pm Operator: McLean
Sample : Inst : Instrumen
Misc : Multiplr: 1.00
MS Integration Params: BENZO.P
Quant Time: Dec 12 11:56 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