

User: Fakhouri, Morgan
 Westchester Dept. of Labs & Research
 Date: 01/16/2002 Time: 14:37:37

Sample: AD28172
 Analysis: \$B1525 Organic Chemicals - 525.2; lab blank
 Units: ug
 Started: 1/16/02 14:32 Ended: 1/16/02 14:32 Analyst: MF
 Result source: manual entry
 Page: 1

	Component Name	Viol	Result	MDL
1	Hexachlorocyclopentadiene		< MDL	0.10
2	Hexachlorobenzene		< MDL	0.10
3	Simazine		< MDL	0.40
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

Acq On : 17 Dec 2001 4:16 pm

Sample :

Misc :

MS Integration Params: BENZO.P

Quant Time: Dec 17 16:53:23 2001

Vial: 27

Operator: McLean

Inst : Instrumen

Multiplr: 1.00

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

Target Compounds	R.T.	QIon	Response	Conc	Units	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-Dinitrotoluene	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

1127LBA.D 011402.M Tue Jan 15 12:47:16 2002

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\011402.M (RTE Integrator)

Title : Quantitation For Method 525.2

Last Update : Tue Dec 18 09:35:41 2001

Response via : Initial Calibration

