

User: Fakhouri, Morgan
Westchester Dept. of Labs & Research
Date: 01/17/2002 Time: 14:02:12

Sample: AD28316
Analysis: \$B1525 Organic Chemicals - 525.2; lab blank
Units: ug
Started: 1/17/02 13:59 Ended: 1/17/02 13:59 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		0.40
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		4.60		
20	Surr_Triphenyl phosphate		5.28		
21	Surr_Perylene-d12		5.35		

Quantitation Report (QT Reviewed)

Data File : C:\MSDCHEM\1\DATA\121701\1127LBB.D

Vial: 43

Acq On : 18 Dec 2001 8:03 am

Operator: McLean

Sample : 11/27

Inst : Instrumen

Misc :

Multiplr: 1.00

MS Integration Params: BENZO.P

Quant Time: Dec 18 08:40:22 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	6782380	5.00	ug/L	0.00
9) Phenanthrene D-10	22.53	188	10470924	5.00	ug/L	0.00
20) Chrysene D-12	30.31	240	7399684	5.00	ug/L	0.00

System Monitoring Compounds

2) 1,3-Dimethyl-2-nitrobenzen	13.13	134	1594209	4.60	ug/L	0.00
Spiked Amount	5.000		Recovery	=	92.00%	
19) Triphenyl Phosphate surr	29.59	326	1682675	5.28	ug/L	0.00
Spiked Amount	5.000		Recovery	=	105.60%	
21) Perylene D-12 surr	34.32	264	5233721	5.35	ug/L	0.00
Spiked Amount	5.000		Recovery	=	107.00%	

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.	d	
5) EPTC	0.00	128	0	N.D.		
6) 2,6-Dinitrotoluene	0.00	165	0	N.D.	d	
7) 2,4-Dinitrotoluene	0.00	165	0	N.D.	d	
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	29.52	129	10927	0.39	ug/L	# 10
23) Bis (2-Ethylhexyl) phthala	30.96	149	325916	0.52	ug/L	94
24) Benzo (a) pyrene	0.00	252	0	N.D.	d	

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

1127LBB.D 011402.M Thu Jan 17 14:03:06 2002

Quantitation Report (QT Reviewed)

Data File : C:\MSDCHEM\1\DATA\121701\1127LBB.D

Vial: 43

Acq On : 18 Dec 2001 8:03 am

Operator: McLean

Sample : 11/27

Inst : Instrumen

Misc :

Multiplr: 1.00

MS Integration Params: BENZO.P

Quant Time: Jan 15 14:27 2002

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

