User: Bohlk, Ted
Westchester Dept. of Labs & Research
Date: 01/10/2002 Time: 09:56:57

Sample: AD27491
Analysis: SB1525 Organic Chemicals - 525.2; lab blank
Units: ug
Started: 11/19/01 09:52 Ended: 12/6/01 09:52 Analyst: TB
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-nitrobenze		4.0		
20	Surr_Triphenyl phosphate		4.9		
21	Surr_Perylene-d12		4.4		

Quantitation Report (QT Reviewed)

Data File : C:\MSDCHEM\1\DATA\120601\1119LB.D Vial: 8
Acq On : 6 Dec 2001 9:15 pm Operator: McLean
Sample : lb Inst : Instrumen
Misc : Multiplr: 1.00
MS Integration Params: BENZO.P
Quant Time: Dec 06 21:52:46 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	4526595	5.00	ug/L	0.00
9) Phenanthrene D-10	22.65	188	8071121	5.00	ug/L	0.00
20) Chrysene D-12	30.43	240	5181540	5.00	ug/L	0.00
System Monitoring Compounds						
2) 1,3-Dimethyl-2-nitrobenzen	13.23	134	954917	3.97	ug/L	0.00
Spiked Amount 5.000			Recovery =	79.40%		
19) Triphenyl Phosphate surr	29.70	326	1381912	4.94	ug/L	0.00
Spiked Amount 5.000			Recovery =	98.80%		
21) Perylene D-12 surr	34.45	264	3213658	4.42	ug/L	0.00
Spiked Amount 5.000			Recovery =	88.40%		
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	17.93	165	21010	0.35	ug/L #	47
7) 2,4-Dinitotoluene	0.00	165	0	N.D. d		
8) Molinate	19.09	126	5102	0.01	ug/L #	1
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	26.76	160	36238	0.11	ug/L #	1
18) 4,4'-DDE	0.00	246	0	N.D. d		
22) Bis (2-Ethylhexyl) adipate	29.62	129	10794	0.12	ug/L #	1
23) Bis (2-Ethylhexyl) phthala	31.07	149	320331	0.41	ug/L	98
24) Benzo (a) pyrene	0.00	252	0	N.D. d		

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

Data File : C:\MSDCHEM\1\DATA\120601\1119LB.D Vial: 8
Acq On : 6 Dec 2001 9:15 pm Operator: McLean
Sample : 1b Inst : Instrumen
Misc : Multiplr: 1.00
MS Integration Params: BENZO.P
Quant Time: Jan 10 8:56 2002 Quant Results File: 120501.RES

Method : C:\MSDCHEM\1\5973N\5080102.M (RTE Integrator)
Title : Quantitation For Method 508gcscan
Last Update : Wed Jan 09 14:35:53 2002
Response via : Initial Calibration