

User: Fakhouri, Morgan
Westchester Dept. of Labs & Research
Date: 01/29/2002 Time: 10:01:58

Sample: AD28105
Analysis: \$B1525 Organic Chemicals - 525.2; lab blank
Units: ug
Started: 1/29/02 09:59 Ended: 1/29/02 09:59 Analyst: MF
Result source: manual entry
Page: 1

Original data

#	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		4.14		
20	Surr_Triphenyl phosphate		5.45		
21	Surr_Perylene-d12		4.08		

Quantitation Report (QT Reviewed)

Data File : C:\MSDCHEM\1\DATA\121001B\1121LB.D

Acq On : 10 Dec 2001 8:10 pm

Sample : 1b

Misc :

MS Integration Params: BENZO.P

Quant Time: Dec 10 20:47:17 2001

Vial: 6

Operator: McLean

Inst : Instrumen

Multiplr: 1.00

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	6235173	5.00	ug/L	0.00
9) Phenanthrene D-10	22.66	188	8074956	5.00	ug/L	0.02
20) Chrysene D-12	30.44	240	6127521	5.00	ug/L	0.02

System Monitoring Compounds

2) 1,3-Dimethyl-2-nitrobenzen	13.24	134	1369625	4.14	ug/L	0.02
Spiked Amount	5.000		Recovery	=	82.80%	
19) Triphenyl Phosphate surr	29.71	326	1525056	5.45	ug/L	0.02
Spiked Amount	5.000		Recovery	=	109.00%	
21) Perylene D-12 surr	34.47	264	3507922	4.08	ug/L	0.03
Spiked Amount	5.000		Recovery	=	81.60%	

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-Dinitrotoluene	19.08	165	17060	0.34 ug/L #	41
8) Molinate	19.08	126	13052	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	26.77	160	46471	0.14 ug/L #	8
18) 4,4'-DDE	0.00	246	0	N.D. d	
22) Bis (2-Ethylhexyl) adipate	29.63	129	6567	0.12 ug/L #	1
23) Bis (2-Ethylhexyl) phthala	31.08	149	70402	0.21 ug/L	96
24) Benzo (a) pyrene	0.00	252	0	N.D. d	

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

1121LB.D 012402B.M Mon Jan 28 15:06:57 2002

Quantitation Report (QT Reviewed)

Data File : C:\MSDCHEM\1\DATA\121001B\1121LB.D

Vial: 6

Acq On : 10 Dec 2001 8:10 pm

Operator: McLean

Sample : lb

Inst : Instrumen

Misc :

Multiplr: 1.00

MS Integration Params: BENZO.P

Quant Time: Jan 28 15:06 2002

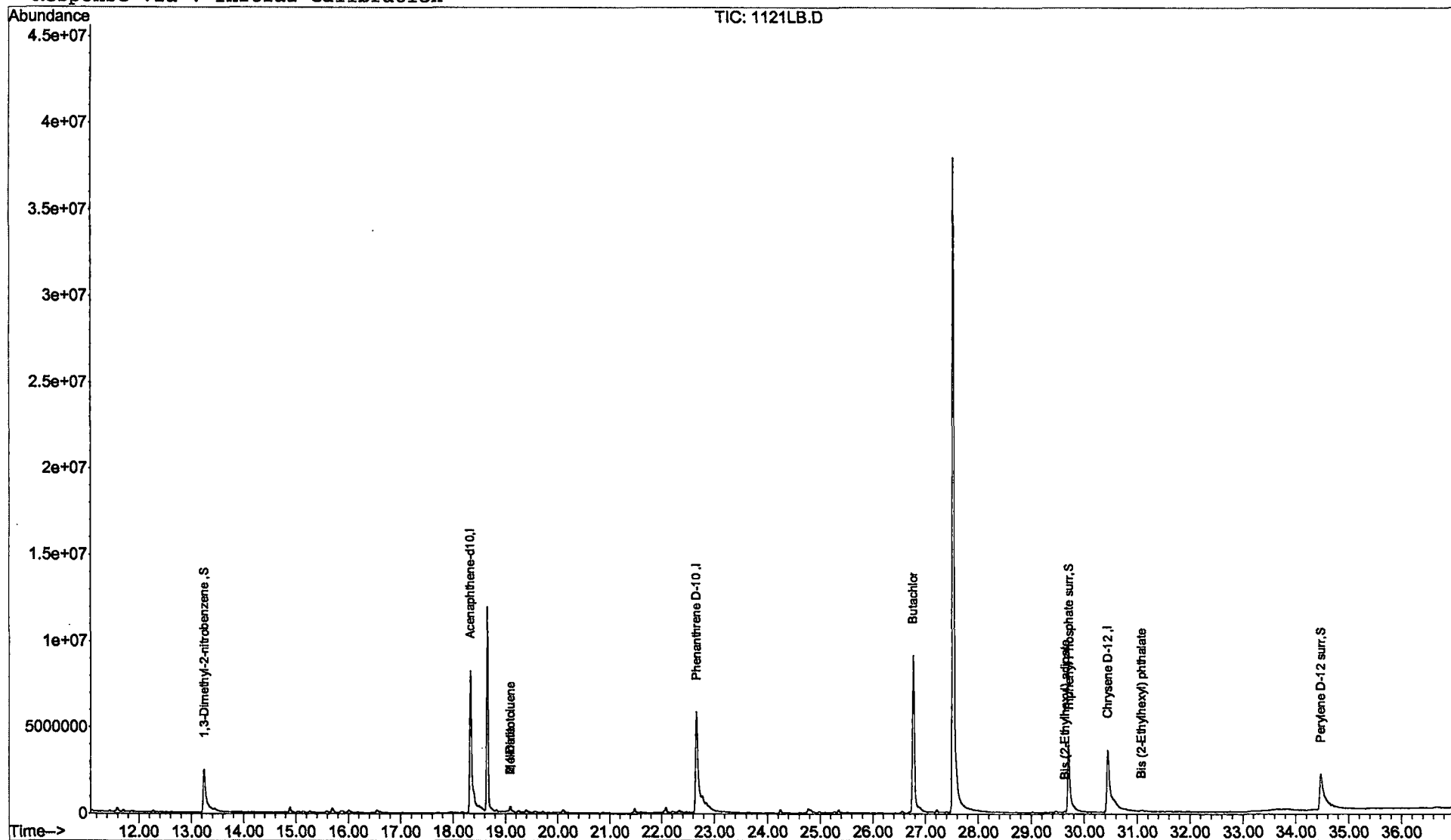
Quant Results File: 120501.RES

Method : C:\MSDCHEM\1\5973N\012402B.M (RTE Integrator)

Title : Quantitation For Method 525.2

Last Update : Fri Jan 25 11:19:14 2002

Response via : Initial Calibration



Quantitation Report (QT Reviewed)

Data File : C:\MSDCHEM\1\DATA\012502A\LB1121R.D

Acq On : 25 Jan 2002 2:39 pm

Sample :

Misc :

MS Integration Params: rteint.p

Quant Time: Jan 28 11:43:32 2002

Vial: 1

Operator: McLean

Inst : Instrumen

Multiplr: 1.00

Quant Results File: 012402B.RES

Reinjection data

Quant Method : C:\MSDCHEM\1\5973N\012402B.M (RTE Integrator)

Title : Quantitation For Method 525.2

Last Update : Fri Jan 25 11:19:14 2002

Response via : Initial Calibration

DataAcq Meth : 012402DF

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev (Min)
1) Acenaphthene-d10	17.97	164	15473729	5.00	ug/L	0.00
9) Phenanthrene D-10	22.27	188	20346992	5.00	ug/L	0.00
20) Chrysene D-12	30.03	240	15149429	5.00	ug/L	0.00

System Monitoring Compounds

2) 1,3-Dimethyl-2-nitrobenzen	12.89	134	4018547	5.07	ug/L	0.00
Spiked Amount	5.000		Recovery	=	101.40%	
19) Triphenyl Phosphate	29.33	326	3137880	5.12	ug/L	0.00
Spiked Amount	5.000		Recovery	=	102.40%	
21) Perylene D-12	34.03	264	10064447	4.68	ug/L	0.00
Spiked Amount	5.000		Recovery	=	93.60%	

Target Compounds

	R.T.	QIon	Response	Conc	Units	Qvalue
3) Hexachlorocyclopentadiene	0.00	237	0	N.D.	d	
4) Hexachlorobenzene	21.14	284	6968	N.D.		
5) EPTC	0.00	128	0	N.D.		
6) 2,6-Dinitrotoluene	17.60	165	31524	0.86	ug/L	30
7) 2,4-Dinitrotoluene	18.70	165	6944	0.91	ug/L	21
8) Molinate	18.90	126	9230	Below Cal	#	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.	d	
17) Butachlor	26.38	160	99688	0.46	ug/L	90
18) 4,4'-DDE	0.00	246	0	N.D.	d	
22) Bis (2-Ethylhexyl) adipate	29.33	129	46679	0.52	ug/L	33
23) Bis (2-Ethylhexyl) phthala	30.69	149	95213	0.40	ug/L	60
24) Benzo (a) pyrene	0.00	252	0	N.D.	d	

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

LB1121R.D 012402B.M

Mon Jan 28 11:45:24 2002

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Quantitation Report (QT Reviewed)

Data File : C:\MSDCHEM\1\DATA\012502A\LB1121R.D
Acq On : 25 Jan 2002 2:39 pm
Sample :
Misc :
MS Integration Params: rteint.p
Quant Time: Jan 28 11:45 2002

Vial: 1
Operator: McLean
Inst : Instrumen
Multiplr: 1.00

Quant Results File: 012402B.RES

Method : C:\MSDCHEM\1\5973N\012402B.M (RTE Integrator)
Title : Quantitation For Method 525.2
Last Update : Fri Jan 25 11:19:14 2002
Response via : Initial Calibration

