User: Bohlk, Ted
Westchester Dept. of Labs & Research
Date: 01/10/2002 Time: 09:00:03

Sample: AD27491
Analysis: S525 Organic Chemicals - 525.2
Units: ug
Started: 11/19/01 08:54 Ended: 12/6/01 08:54 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		0.83 — 0.4		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		4.2		1.0
20	Surr_Triphenyl phosphate		6.2		1.0
21	Surr_Perylene-d12		4.9		1.0

Quantitation Report (QT Reviewed)

Data File : C:\MSDCHEM\1\DATA\120601\27491.D Vial: 21
Acq On : 7 Dec 2001 7:40 am Operator: McLean
Sample : 11/19 sample Inst : Instrumen
Misc : Multiplr: 1.00
MS Integration Params: BENZO.P
Quant Time: Dec 07 08:17:40 2001 Quant Results File: 120501.RES

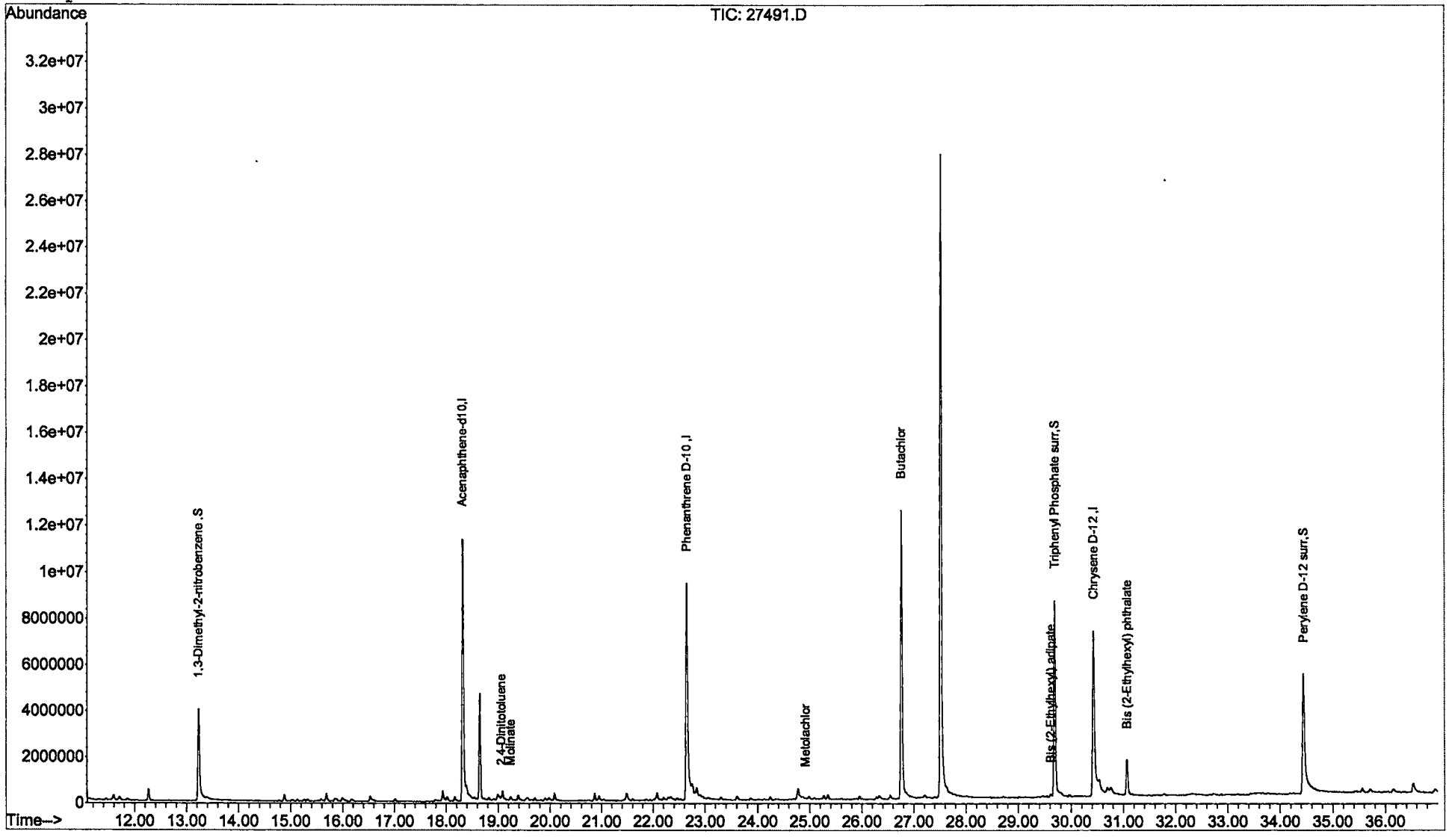
Quant Method : C:\MSDCHEM\1\5973N\120501.M (RTE Integrator)
Title : Quantitation For Method 525.2
Last Update : Thu Dec 06 10:31:04 2001
Response via : Initial Calibration
DataAcq Meth : 120501

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Acenaphthene-d10	18.32	164	6573028	5.00	ug/L	0.00
9) Phenanthrene D-10	22.64	188	10146860	5.00	ug/L	0.00
20) Chrysene D-12	30.42	240	7248552	5.00	ug/L	0.00
System Monitoring Compounds						
2) 1,3-Dimethyl-2-nitrobenzen	13.23	134	1450053	4.16	ug/L	0.00
Spiked Amount 5.000			Recovery = 83.20%			
19) Triphenyl Phosphate surr	29.70	326	2166336	6.16	ug/L	0.01
Spiked Amount 5.000			Recovery = 123.20%			
21) Perylene D-12 surr	34.44	264	4976311	4.89	ug/L	0.00
Spiked Amount 5.000			Recovery = 97.80%			
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	19.05	165	14753	0.33	ug/L #	42
8) Molinate	19.23	126	5315	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	24.93	162	7244	0.01	ug/L #	38
17) Butachlor	26.76	160	46854	0.11	ug/L #	9
18) 4,4'-DDE	0.00	246	0	N.D. d		
22) Bis (2-Ethylhexyl) adipate	29.62	129	28070	0.13	ug/L #	28
23) Bis (2-Ethylhexyl) phthala	31.07	149	1224308	0.83	ug/L	99
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\27491.D Vial: 21
Acq On : 7 Dec 2001 7:40 am Operator: McLean
Sample : 11/19 sample Inst : Instrumen
Misc : Multiplr: 1.00
MS Integration Params: BENZO.P
Quant Time: Jan 10 8:55 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



Comments for Sample AD27491 - Analysis $525

Westchester Dept. of Labs & Research
User: Vinci, David L
Date: 01-10-2002 Time: 12:28:50

The recoveries for 6 of the 18 compounds in the LCS Standard were outside of their acceptance ranges.