User: Fakhouri, Morgan
Westchester Dept. of Labs & Research
Date: 01/16/2002 Time: 14:29:28

Sample: AD28172
Analysis: $525 Organic Chemicals - 525.2
Units: ug
Started: 1/16/02 14:22 Ended: 1/16/02 14:22 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		4.95	1.0
20	Surr_Triphenyl phosphate		5.88	1.0
21	Surr_Perylene-d12		5.28	1.0

Quantitation Report (Not Reviewed)

Data File : C:\MSDCHEM\1\DATA\121701\28172.D Vial: 34
Acq On : 17 Dec 2001 9:31 pm Operator: McLean
Sample : 1126 Inst : Instrumen
Misc : Multiplr: 1.00
MS Integration Params: BENZO.P
Quant Time: Dec 17 22:08:46 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

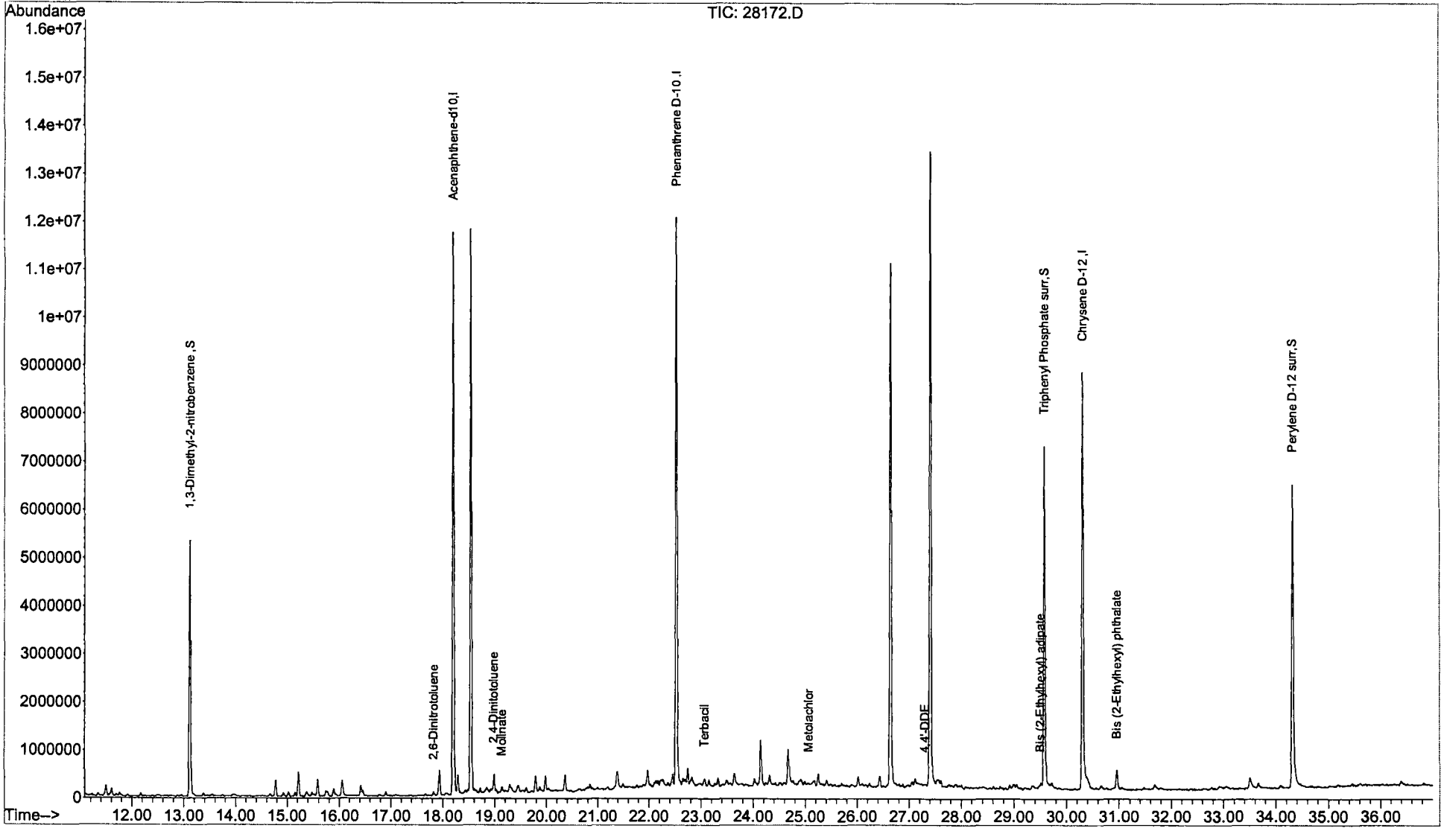
Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Acenaphthene-d10	18.20	164	5556162	5.00	ug/L	0.00
9) Phenanthrene D-10	22.53	188	9145596	5.00	ug/L	0.00
20) Chrysene D-12	30.30	240	6005469	5.00	ug/L	0.00
System Monitoring Compounds						
2) 1,3-Dimethyl-2-nitrobenzen	13.13	134	1406848	4.95	ug/L	0.00
Spiked Amount 5.000			Recovery = 99.00%			
19) Triphenyl Phosphate surr	29.58	326	1636257	5.88	ug/L	0.00
Spiked Amount 5.000			Recovery = 117.60%			
21) Perylene D-12 surr	34.31	264	4194825	5.28	ug/L	0.00
Spiked Amount 5.000			Recovery = 105.60%			
Target Compounds						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	17.82	165	16674	0.36	ug/L #	39
7) 2,4-Dinitotoluene	18.97	165	3254	0.32	ug/L #	21
8) Molinate	19.14	126	5775	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	23.07	161	17044	0.06	ug/L #	27
14) Acetochlor	0.00	146	0	N.D.		
15) Metribuzin	0.00	198	0	N.D. d		
16) Metolachlor	25.06	162	7291	0.01	ug/L #	38
17) Butachlor	0.00	160	0	N.D. d		
18) 4,4'-DDE	27.30	246	5209	0.01	ug/L #	23
22) Bis (2-Ethylhexyl) adipate	29.52	129	25335	0.41	ug/L #	21
23) Bis (2-Ethylhexyl) phthala	30.96	149	340086	0.55	ug/L	97
24) Benzo (a) pyrene	0.00	252	0	N.D. d		

(#) = qualifier out of range (m) = manual integration
28172.D 011402.M Wed Jan 16 14:24:23 2002 Page 1

Quantitation Report (Not Reviewed)

Data File : C:\MSDCHEM\1\DATA\121701\28172.D Vial: 34
Acq On : 17 Dec 2001 9:31 pm Operator: McLean
Sample : 1126 Inst : Instrumen
Misc : Multiplr: 1.00
MS Integration Params: BENZO.P
Quant Time: Jan 15 12:55 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



Comments for Sample AD28172 - Analysis $525

Westchester Dept. of Labs & Research
User: Vinci, David L
Date: 01-22-2002 Time: 08:37:57

The breakdown for Endrin in the Breakdown Standard was 21%. The recoveries for 3 of the 18 compounds in the LCS Standard were outside of their acceptance ranges.