User: Fakhouri, Morgan
Westchester Dept. of Labs & Research
Date: 12/11/2001 Time: 11:17:09

Sample: AD27547
Analysis: $C1525 Organic Chemicals - 525.2; cal. std.
Units: ug
Started: 11/16/01 10:58 Ended: 12/5/01 10:58 Analyst: MF
Page: 1

CCV

	Component Name	Viol	Result	Qualifier	MDL
1	Hexachlorocyclopentadiene		6.63		0.10
2	Hexachlorobenzene		4.96		0.10
3	Simazine		5.87		0.07
4	Atrazine		6.23		0.10
5	Alachlor		6.41		0.20
6	bis(2-Ethylhexyl)adipate		6.62		0.60
7	bis(2-Ethylhexyl)phthalate		~~6.86~~ 6.24 DW 12/11/1.		0.60
8	Benzo(a)pyrene		5.66		0.20
9	EPTC		5.32		1.0
10	2,6-Dinitrotoluene		4.98		2.0
11	2,4-Dinitrotoluene		5.32		2.0
12	Molinate		5.47		0.90
13	Terbacil		5.66		2.0
14	Acetochlor		6.41		2.0
15	Metribuzin		6.83		0.15
16	Metolachlor		6.73		0.75
17	Butachlor		5.29		0.40
18	4,4-DDE		5.98		0.80
19	Surr_1,3-Dimethyl-2-nitrobenzen		3.71		
20	Surr_Triphenyl phosphate		5.40		
21	Surr_Perylene-d12		7.54		

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Quantitation Report (QT Reviewed)

Data File : C:\MSDCHEM\1\DATA\120501\CAL6.D Vial: 4
Acq On : 5 Dec 2001 5:38 pm Operator: McLean
Sample : 5.0 cc Inst : Instrumen
Misc : Multiplr: 1.00
MS Integration Params: BENZO.P
Quant Time: Dec 11 09:18:20 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	18.31	164	6382601	5.00	ug/L	0.00
9) Phenanthrene D-10	22.64	188	8945263	5.00	ug/L	0.00
20) Chrysene D-12	30.42	240	4551950	5.00	ug/L	0.00
System Monitoring Compounds						
2) 1,3-Dimethyl-2-nitrobenzen	13.23	134	1257886	3.71	ug/L	0.00
Spiked Amount 5.000			Recovery = 74.20%			
19) Triphenyl Phosphate surr	29.69	326	1675568	5.40	ug/L	0.00
Spiked Amount 5.000			Recovery = 108.00%			
21) Perylene D-12 surr	34.44	264	4818293	7.54	ug/L	0.00
Spiked Amount 5.000			Recovery = 150.80%			
Target Compounds						Qvalue
3) Hexachlorocyclopentadiene	15.81	237	1305324	6.63	ug/L	97
4) Hexachlorobenzene	21.51	284	2370592	4.96	ug/L	97
5) EPTC	16.28	128	3570278	5.32	ug/L	99
6) 2,6-Dinitrotoluene	17.88	165	1244123	4.98	ug/L	97
7) 2,4-Dinitotoluene	19.03	165	1234918	4.20	ug/L	99
8) Molinate	19.20	126	5155440	5.47	ug/L	91
10) Simazine	22.04	201	1471307	5.87	ug/L	98
11) Atrazine	22.16	200	2283166	6.23	ug/L	97
12) Alachlor	24.02	160	2168655	6.41	ug/L	93
13) Terbacil	22.98	161	2163059	5.66	ug/L	100
14) Acetochlor	23.80	146	1496600	6.48	ug/L	99
15) Metribuzin	23.84	198	2666272	6.83	ug/L	95
16) Metolachlor	24.94	162	6784795	6.73	ug/L	99
17) Butachlor	26.77	160	1961924	5.29	ug/L	86
18) 4,4'-DDE	27.40	246	2213634	5.98	ug/L	97
22) Bis (2-Ethylhexyl) adipate	29.62	129	4532620m	6.62	ug/L	
23) Bis (2-Ethylhexyl) phthala	31.07	149	8086961m	6.24	ug/L	
24) Benzo (a) pyrene	34.29	252	5286074m	5.66	ug/L	

3.5-6.5 3/18 Lab

(#) = qualifier out of range (m) = manual integration
CAL6.D 120501.M Tue Dec 11 14:57:50 2001 Page 1

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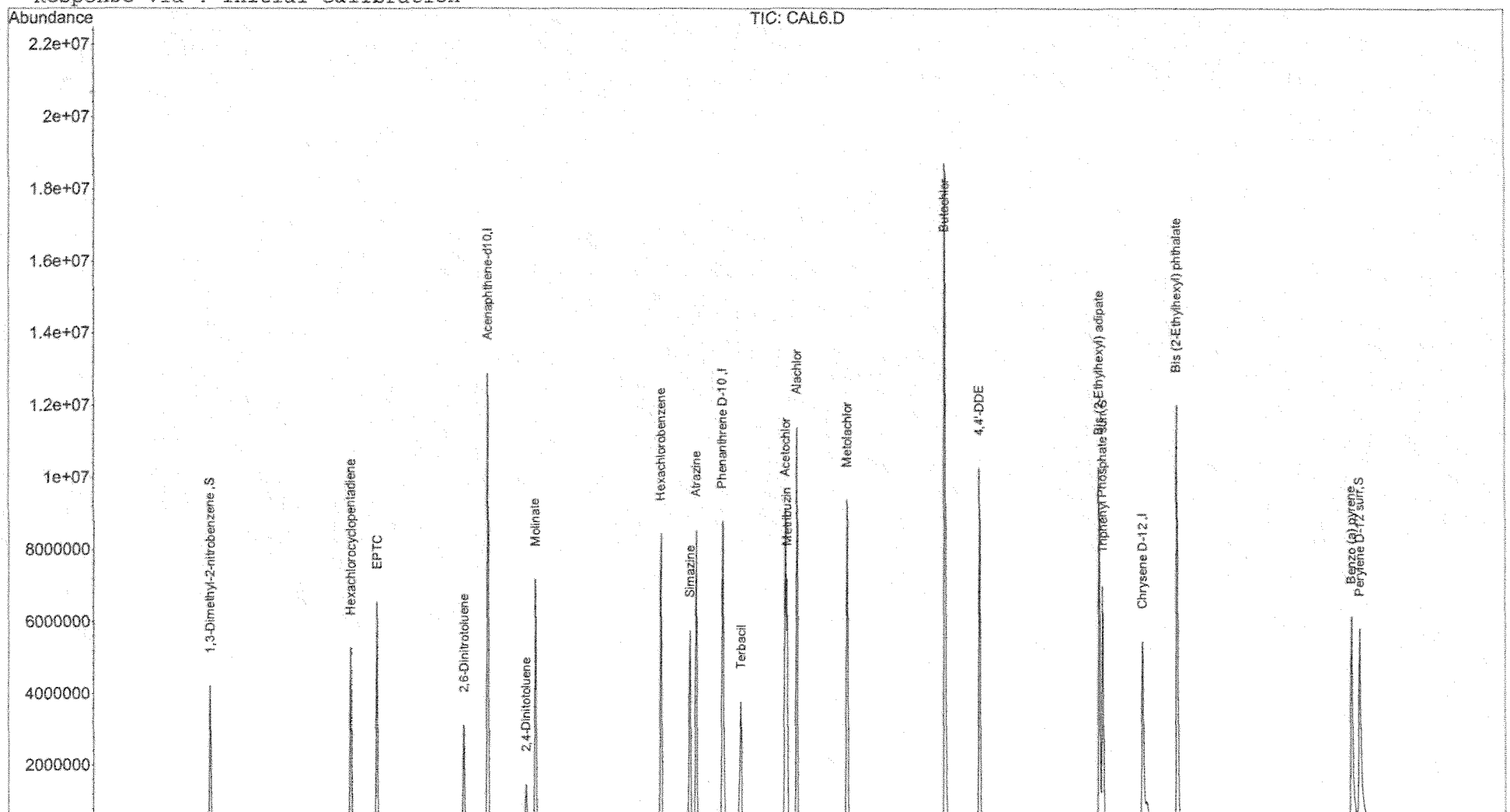
Quantitation Report (QT Reviewed)

Data File : C:\MSDCHEM\1\DATA\120501\CAL6.D Vial: 4
Acq On : 5 Dec 2001 5:38 pm Operator: McLean
Sample : 5.0 cc Inst : Instrumen
Misc : Multiplr: 1.00
MS Integration Params: BENZO.P
Quant Time: Dec 11 14:57 2001 Quant Results File: 120501.RES

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



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NYC-WTC_000057729