

User: Vinci, David L
Westchester Dept. of Labs & Research
Date: 02/27/2002 Time: 14:28:49

Sample: AE00949
Analysis: \$GCMS GCMS Scan For Unknowns
Units: ug
Started: 2/27/02 14:28 Ended: 2/27/02 14:28 Analyst: DLV
Page: 1

	Component Name	Viol	Result	Qualifier	MDL
1	Other unknown compounds		.		
2	Surr_2,4-DDD		70		5.0

Comments for Sample AE00949 - Analysis \$GCMS

Westchester Dept. of Labs & Research

User: Vinci, David L

Date: 02-27-2002 Time: 14:28:55

The GCMS scan, from 35 to 550 amu, does not reveal the presence of unusual compounds. The recovery for 1 of the 43 compounds in the LCS Standard was below its acceptance range.

Data File : C:\MSDCHEM\1\DATA\022202\0949.D

Vial: 16

Acq On : 23 Feb 2002 1:17 am

Operator: McLean

Sample : 1/14/02

Inst : Instrumen

Misc :

Multiplr: 1.00

MS Integration Params: BENZO.P

Quant Time: Feb 27 06:26:28 2002

Quant Results File: 5080202.RES

Quant Method : C:\MSDCHEM\1\5973N\5080202.M (RTE Integrator)

Title : Quantitation For Method 508gcscan

Last Update : Mon Feb 25 19:17:14 2002

Response via : Initial Calibration

DataAcq Meth : 508SCAN

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	9.58	150	942770	20.00	ug/L	0.00
10) Naphthalene-d8	13.05	136	2505961	20.00	ug/L	0.00
17) Acenaphthene-d10	18.16	164	1347792	20.00	ug/L	0.00
29) Phenanthrene-d10	22.52	188	2073258	20.00	ug/L	0.00
38) Chrysene-d12	30.33	240	1934777	20.00	ug/L	-0.03
44) Perylene-d12	34.21	264	1310726	20.00	ug/L	0.00

System Monitoring Compounds

37) 2,4-ddd surr	27.47	235	115939	3.52	ug/L	0.00
Spiked Amount	5.000		Recovery	=	70.40%	

Target Compounds

					Qvalue
2) N-nitroso-dimethyl amine	0.00	74	0	N.D.	
3) bis(2-chloroethyl) ether	0.00	93	0	N.D.	
4) 1,3 - Dichlorobenze	0.00	146	0	N.D.	
5) 1,4 - Dichlorobenzene	0.00	146	0	N.D.	
6) 1,2 - Dichlorobenzene	0.00	146	0	N.D.	
7) 2,2' - oxybis (1-Chloropr	0.00	45	0	N.D.	
8) N-Nitroso-di-n-propylamine	0.00	70	0	N.D.	
9) Hexachloroethane	0.00	117	0	N.D.	
11) Nitrobenzene	0.00	77	0	N.D.	
12) Isophorone	0.00	82	0	N.D.	
13) bis(2-Chloroethoxy) methan	0.00	93	0	N.D.	
14) 1,2,4-Trichlorobenzene	0.00	180	0	N.D.	
15) Naphthalene	0.00	128	0	N.D.	
16) Hexachlorobutadiene	0.00	225	0	N.D.	
18) Hexachlorocyclopentadiene	0.00	237	0	N.D.	
19) 2-chloronaphthalene	0.00	162	0	N.D.	
20) Dimethylphthalate	0.00	163	0	N.D.	
21) 2,6-Dinitrotoluene	18.16	165	170040	9.07 ug/L #	25
22) Acenaphthylene	0.00	152	0	N.D.	
23) Acenaphthene	0.00	153	0	N.D.	
24) 2,4-Dinitrotoluene	0.00	165	0	N.D.	
25) Diethylphthalate	0.00	149	0	N.D.	
26) 4-Chlorophenyl-phenyl ethe	0.00	204	0	N.D.	
27) Fluorene	0.00	166	0	N.D.	
28) Azobenzen/ 1,2-Diphenylhyd	0.00	77	0	N.D.	
30) N-Nitrosodiphenylamine	0.00	169	0	N.D.	
31) 4-Bromophenyl-phenyl ether	0.00	248	0	N.D.	
32) Hexachlorobenzene	0.00	284	0	N.D.	
33) Phenanthrene	0.00	178	0	N.D.	
34) Anthracene	0.00	178	0	N.D.	
35) Di-n-butylphthalate	24.67	149	15180	0.11 ug/L	79
36) Fluoranthene	0.00	202	0	N.D.	
39) Pyrene	0.00	202	0	N.D.	
40) Butylbenzylphthalate	0.00	149	0	N.D.	
41) Benzo (a) anthracene	30.33	228	5250	0.04 ug/L	66
42) Bis (2-ethylhexyl) phthala	30.95	149	2502	0.03 ug/L	48
43) Chrysene	30.33	228	5250	0.04 ug/L	63
45) Di-n-Octylphthalate	0.00	149	0	N.D.	
46) Benzo (b) fluoranthene	0.00	252	0	N.D.	

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

0949.D 5080202.M Wed Feb 27 06:26:35 2002

Data File : C:\MSDCHEM\1\DATA\022202\0949.D

Vial: 16

Acq On : 23 Feb 2002 1:17 am

Operator: McLean

Sample : 1/14/02

Inst : Instrumen

Misc :

Multiplr: 1.00

MS Integration Params: BENZO.P

Quant Time: Feb 27 06:26:28 2002

Quant Results File: 5080202.RES

Quant Method : C:\MSDCHEM\1\5973N\5080202.M (RTE Integrator)

Title : Quantitation For Method 508gcscan

Last Update : Mon Feb 25 19:17:14 2002

Response via : Initial Calibration

DataAcq Meth : 508SCAN

Compound	R.T.	QIon	Response	Conc	Unit	Qvalue
47) Benzo (k) fluoranthene	0.00	252	0	N.D.		
48) Benzo (a) pyrene	34.21	252	5169	0.05 ug/L		64
49) Indeno (1,2,3-cd) pyrene	0.00	276	0	N.D.		
50) Dibenz (a,h) anthracene	0.00	278	0	N.D.		
51) Benzo (g,h,i) perylene	0.00	276	0	N.D.		

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

0949.D 5080202.M

Wed Feb 27 06:26:35 2002

Page 2

Data File : C:\MSDCHEM\1\DATA\022202\0949.D

Vial: 16

Acq On : 23 Feb 2002 1:17 am

Operator: McLean

Sample : 1/14/02

Inst : Instrumen

Misc :

Multiplr: 1.00

MS Integration Params: BENZO.P

Quant Time: Feb 27 6:26 2002

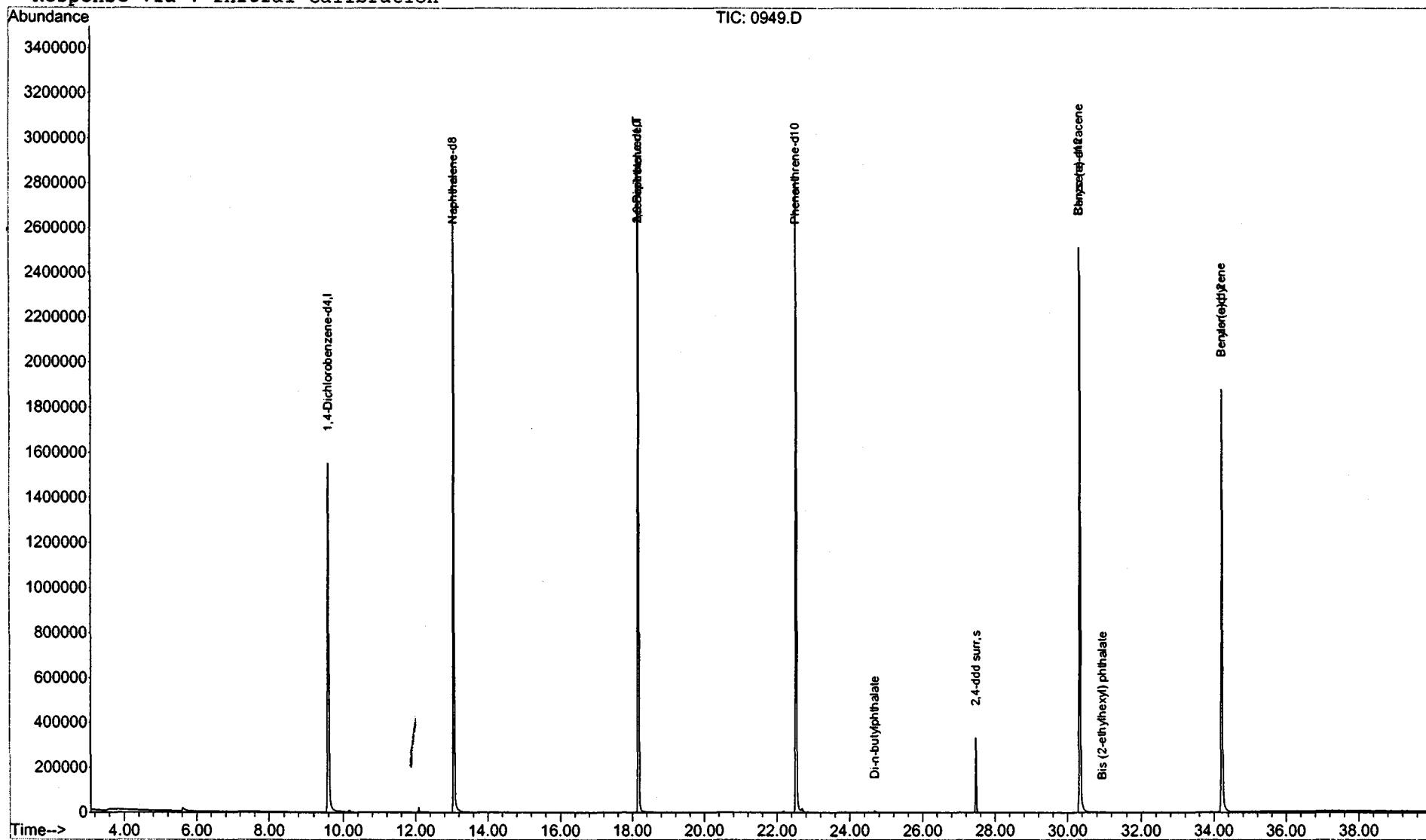
Quant Results File: 5080202.RES

Method : C:\MSDCHEM\1\5973N\5080202.M (RTE Integrator)

Title : Quantitation For Method 508gcscan

Last Update : Mon Feb 25 19:17:14 2002

Response via : Initial Calibration



LSC Area Percent Report

Data File : C:\MSDCHEM\1\DATA\022202\0949.D

Vial: 16

Acq On : 23 Feb 2002 1:17 am

Operator: McLean

Sample : 1/14/02

Inst : Instrumen

Misc :

Multiplr: 1.00

MS Integration Params: LSCINT.P

Method : C:\MSDCHEM\1\5973N\5080202.M (RTE Integrator)

Title : Quantitation For Method 508gcscan

Smoothing : OFF

Filtering: 5

Sampling : 1

Min Area: 1 % of largest Peak

Start Thrs: 0.2

Max Peaks: 100

Stop Thrs : 0

Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >

Peak separation: 5

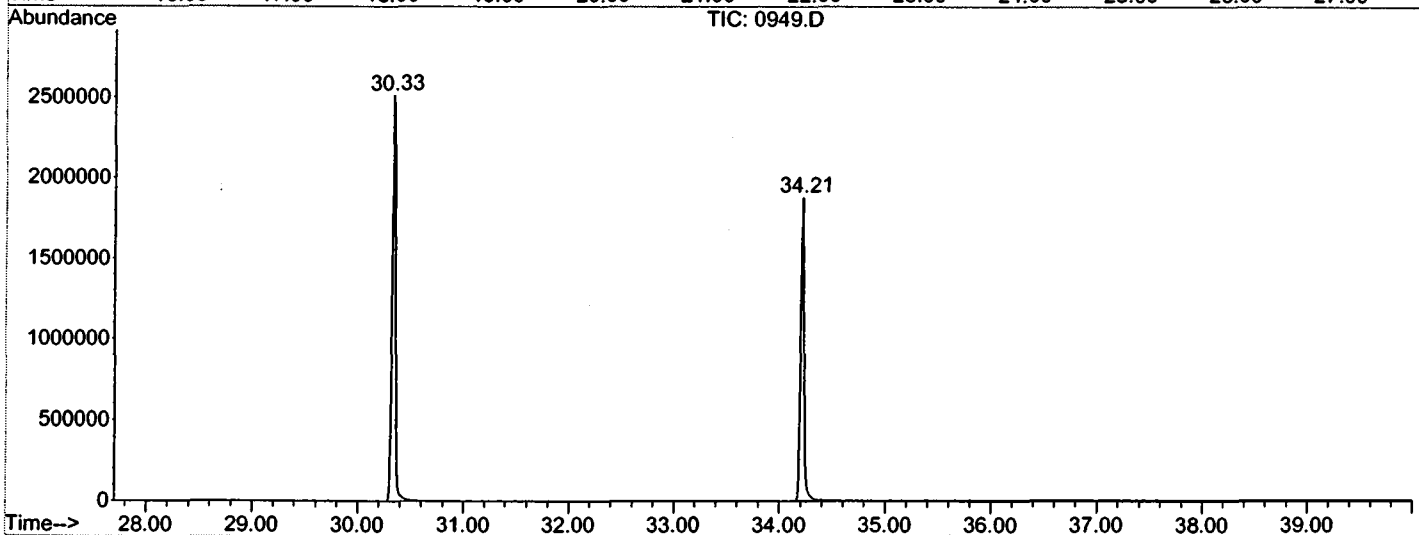
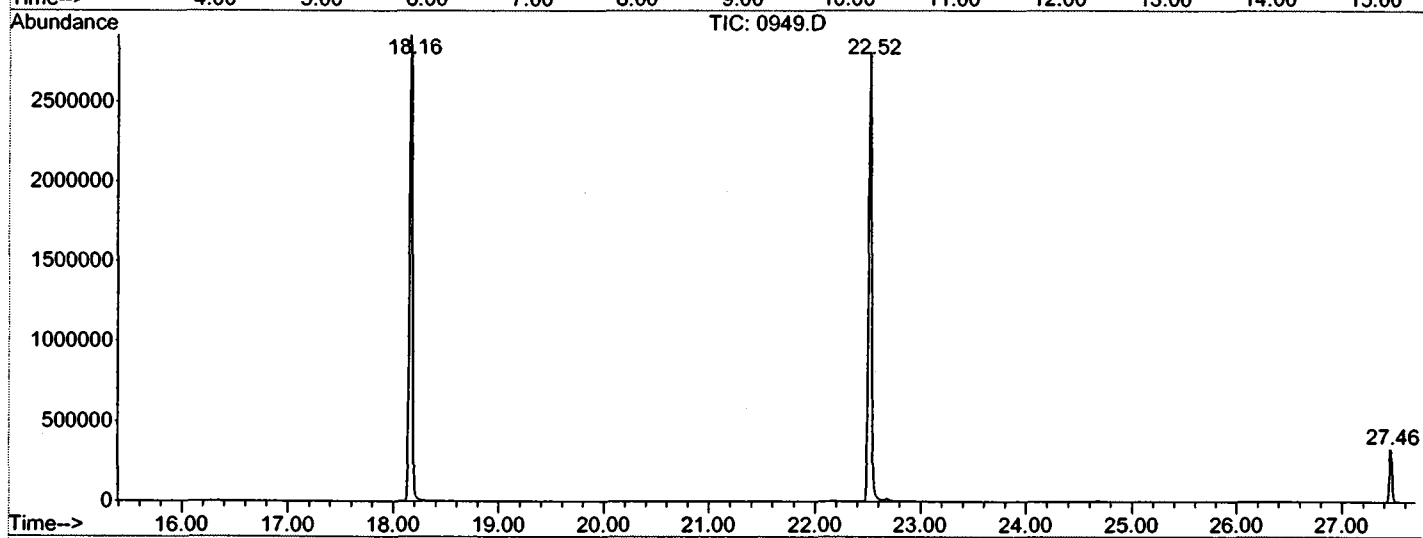
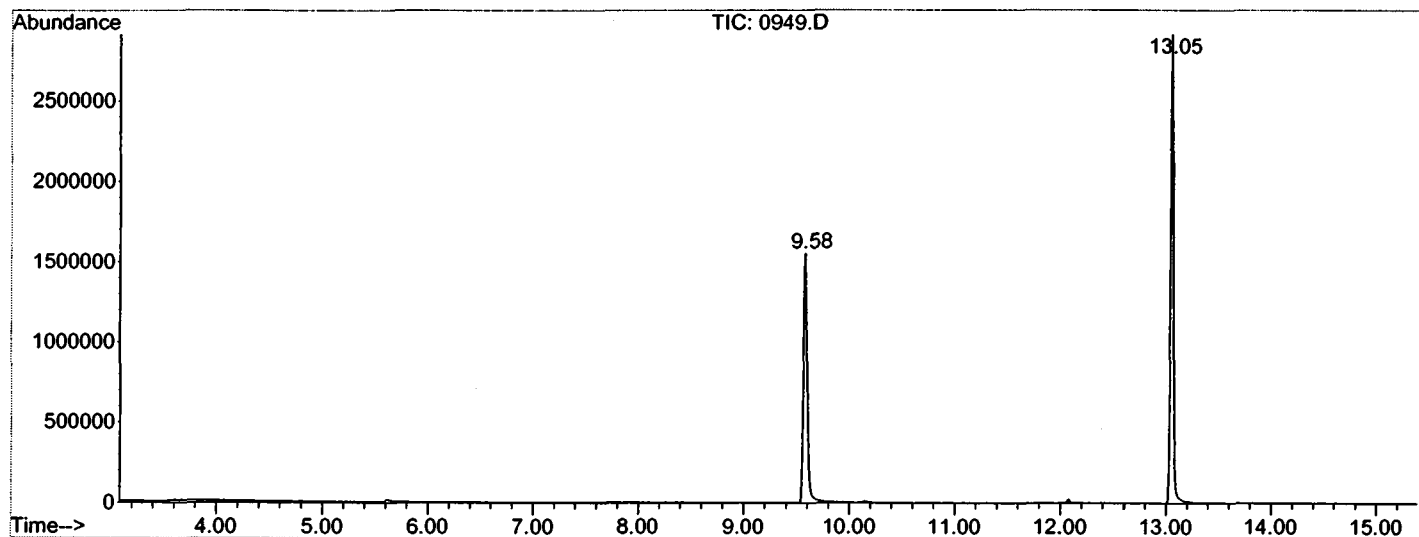
Signal : TIC

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	9.579	712	717	732	rBV	1552821	3913470	68.80%	12.534%
2	13.053	1095	1100	1116	rBV	2912993	5414527	95.19%	17.341%
3	18.160	1658	1663	1674	rBV	2913567	5656540	99.45%	18.116%
4	22.523	2138	2144	2154	rBV2	2808873	5687888	100.00%	18.217%
5	27.457	2684	2688	2695	rBV	330842	599170	10.53%	1.919%
6	30.332	2999	3005	3019	rBV2	2513857	5629370	98.97%	18.029%
7	34.214	3426	3433	3451	rBV	1878896	4322231	75.99%	13.843%

Sum of corrected areas: 31223196

LSC Report - Integrated Chromatogram

File : C:\MSDCHEM\1\DATA\022202\0949.D
 Operator : McLean
 Acquired : 23 Feb 2002 1:17 am using AcqMethod 508SCAN
 Instrument : Instrumen
 Sample Name: 1/14/02
 Misc Info :
 Vial Number: 16
 Quant File :5080202.RES (RTE Integrator)



Tentatively Identified Compound (LSC) summary

Operator ID: McLean Date Acquired: 23 Feb 2002 1:17 am
 Data File: C:\MSDCHEM\1\DATA\022202\0949.D
 Name: 1/14/02
 Misc:
 Method: C:\MSDCHEM\1\5973N\5080202.M (RTE Integrator)
 Title: Quantitation For Method 508gcscan
 Library Searched: C:\DATABASE\WILEY7N.L

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
