

User: Vinci, David L
Westchester Dept. of Labs & Research
Date: 04/05/2002 Time: 15:59:18

Sample: AE06046
Analysis: \$GCMS GCMS Scan For Unknowns
Units: ug
Started: 4/5/02 15:58 Ended: 4/5/02 15:58 Analyst: DLV
Page: 1

	Component Name	Viol	Result	Qualifier	MDL
1	Other unknown compounds				
2	Surr_2,4-DDD		100		5.0

Comments for Sample AE06046 - Analysis \$GCMS

Westchester Dept. of Labs & Research

User: Vinci, David L

Date: 04-05-2002 Time: 15:59:38

The GCMS scan, from 35 to 550 amu, does not reveal the presence of unusual compounds.

Quantitation Report (QT Reviewed)

Data File : C:\MSDCHEM\1\DATA\032202\6046.D

Vial: 17

Acq On : 23 Mar 2002 8:26 am

Operator: McLean

Sample : SAMPLE 6046 3/14/02

Inst : Instrumen

Misc :

Multiplr: 1.00

MS Integration Params: BENZO.P

Quant Time: Apr 05 07:38:34 2002

Quant Results File: 5080302.RES

Quant Method : C:\MSDCHEM\1\5973N\5080302.M (RTE Integrator)

Title : Quantitation For Method 508gcscan

Last Update : Wed Mar 27 06:57:36 2002

Response via : Initial Calibration

DataAcq Meth : 508SCAN

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev (Min)
1) 1,4-Dichlorobenzene-d4	9.55	150	1526172	20.00	ug/L	0.00
10) Naphthalene-d8	13.03	136	4449776	20.00	ug/L	0.00
17) Acenaphthene-d10	18.14	164	2526234	20.00	ug/L	0.00
29) Phenanthrene-d10	22.50	188	3842375	20.00	ug/L	0.00
38) Chrysene-d12	30.31	240	3015817	20.00	ug/L	-0.01
44) Perylene-d12	34.18	264	1576549	20.00	ug/L	0.00

System Monitoring Compounds

37) 2,4-DDD (SURR)	27.43	235	229524	5.01	ug/L	0.00
Spiked Amount	5.000		Recovery	=	100.20%	

Target Compounds

					Qvalue
2) N-nitroso-dimethyl amine	3.26	74	8145	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	0.00	165	0	N.D.	d
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	24.65	149	23841	N.D.	
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	30.31	228	6934	N.D.	
42) Bis (2-ethylhexyl) phthala	30.91	149	10441	N.D.	
43) Chrysene	30.31	228	6934	N.D.	
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

6046.D 5080302.M Fri Apr 05 07:38:46 2002

Quantitation Report (QT Reviewed)

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Operator: McLean

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Quant Method : C:\MSDCHEM\1\5973N\5080302.M (RTE Integrator)

Title : Quantitation For Method 508gcscan

Last Update : Wed Mar 27 06:57:36 2002

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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	34.18	252	5397	N.D.		
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

6046.D 5080302.M

Fri Apr 05 07:38:46 2002

Page 2

Data File : C:\MSDCHEM\1\DATA\032202\6046.D

Vial: 17

Acq On : 23 Mar 2002 8:26 am

Operator: McLean

Sample : SAMPLE 6046 3/14/02

Inst : Instrumen

Misc :

Multiplr: 1.00

MS Integration Params: BENZO.P

Quant Time: Apr 5 7:38 2002

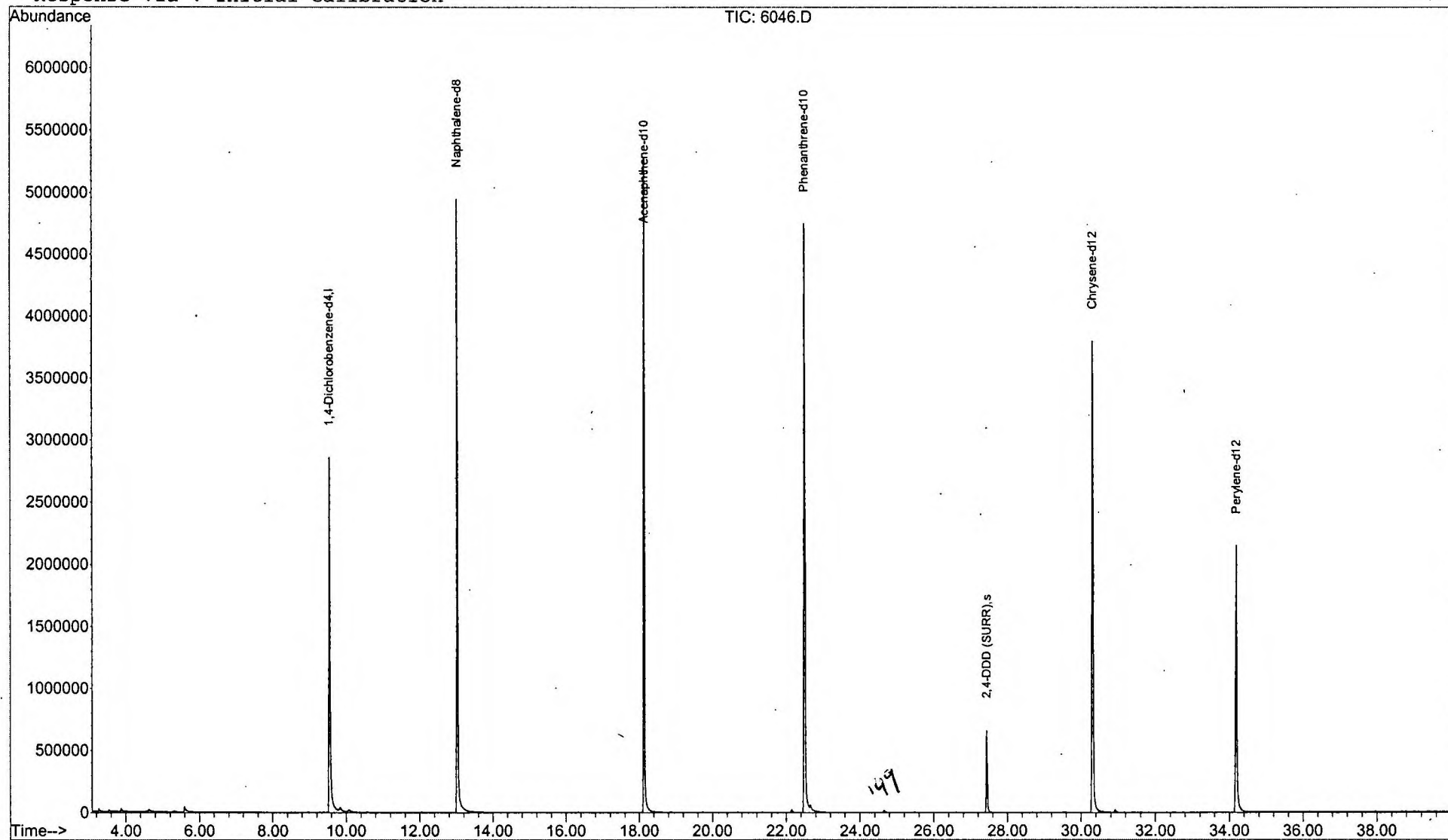
Quant Results File: 5080302.RES

Method : C:\MSDCHEM\1\5973N\5080302.M (RTE Integrator)

Title : Quantitation For Method 508gcscan

Last Update : Wed Mar 27 06:57:36 2002

Response via : Initial Calibration



LSC Area Percent Report

Data File : C:\MSDCHEM\1\DATA\032202\6046.D
Acq On : 23 Mar 2002 8:26 am
Sample : SAMPLE 6046 3/14/02
Misc :
MS Integration Params: LSCINT.P

Vial: 17
Operator: McLean
Inst : Instrumen
Multiplr: 1.00

Method : C:\MSDCHEM\1\5973N\5080302.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

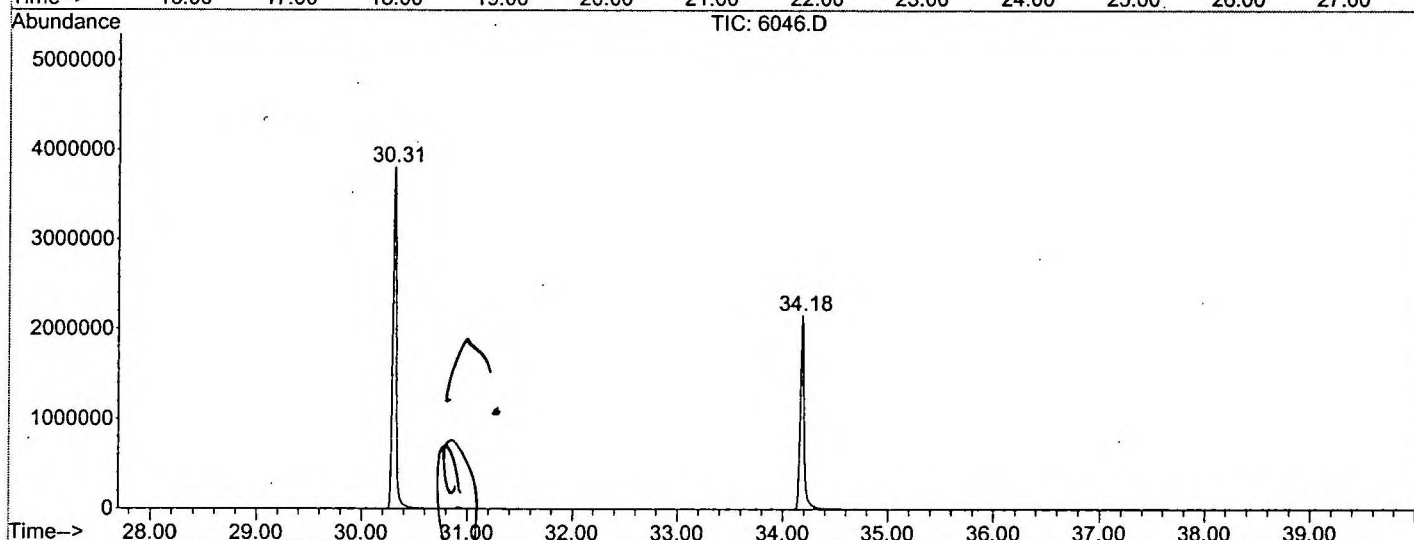
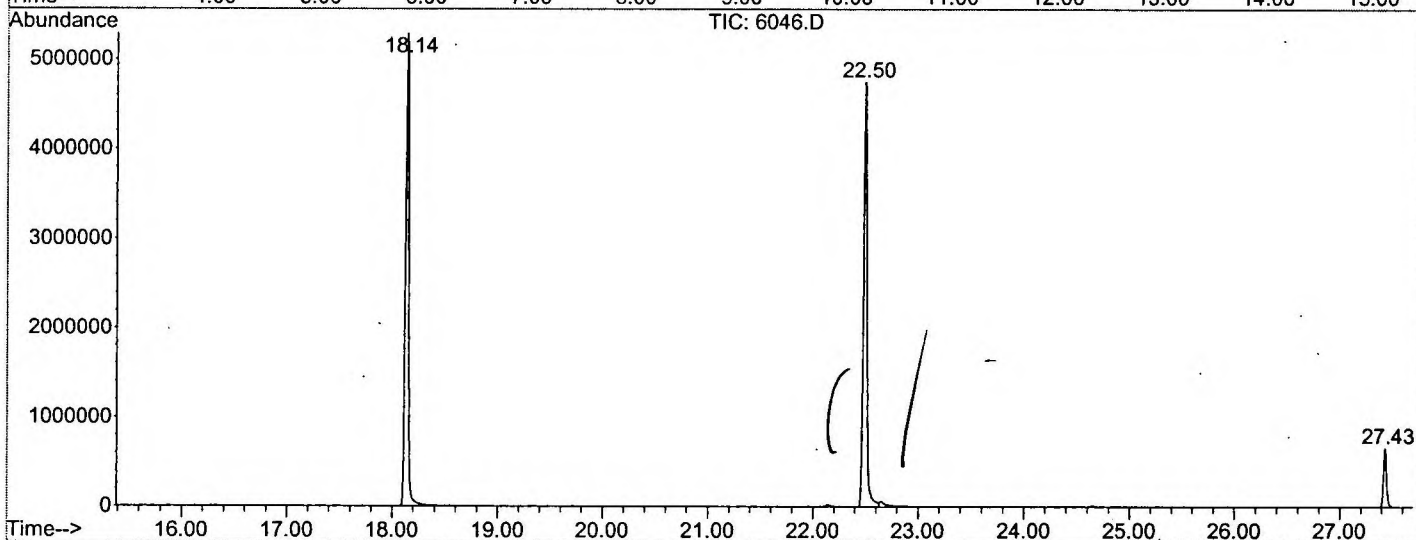
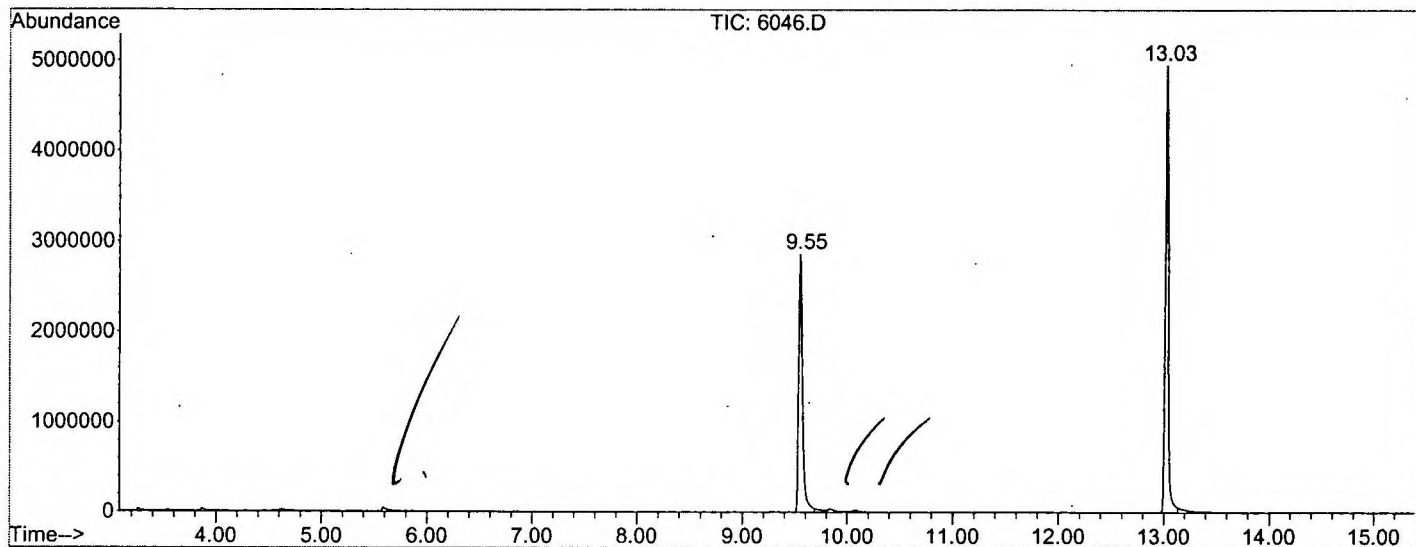
Signal : TIC

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	9.548	1094	1101	1125	rBV	2855704	6955954	63.52%	12.738%
2	13.027	1685	1693	1711	rBV	4947256	10064221	91.90%	18.429%
3	18.138	2553	2563	2583	rBV	5282190	10951398	100.00%	20.054%
4	22.498	3295	3305	3327	rBV2	4748037	10920690	99.72%	19.998%
5	27.433	4137	4145	4162	rBV	656416	1289483	11.77%	2.361%
6	30.313	4623	4635	4661	rBV2	3803281	9086631	82.97%	16.639%
7	34.185	5283	5294	5314	rBV2	2152098	5341010	48.77%	9.780%

Sum of corrected areas: 54609387

LSC Report - Integrated Chromatogram

File : C:\MSDCHEM\1\DATA\032202\6046.D
 Operator : McLean
 Acquired : 23 Mar 2002 8:26 am using AcqMethod 508SCAN
 Instrument : Instrumen
 Sample Name: SAMPLE 6046 3/14/02
 Misc Info :
 Vial Number: 17
 Quant File :5080302.RES (RTE Integrator)



Tentatively Identified Compound (LSC) summary

Operator ID: McLean Date Acquired: 23 Mar 2002 8:26 am
 Data File: C:\MSDCHEM\1\DATA\032202\6046.D
 Name: SAMPLE 6046 3/14/02
 Misc:
 Method: C:\MSDCHEM\1\5973N\5080302.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
