



User: Vinci, David L
 Westchester Dept. of Labs & Research
 Date: 03/21/2002 Time: 15:40:17

Sample: AE05913
 Analysis: \$GCMS GCMS Scan For Unknowns
 Units: ug
 Started: 3/21/02 15:37 Ended: 3/21/02 15:37 Analyst: DLV
 Page: 1

	Component Name	Viol	Result	Qualifier	MDL
1	Other unknown compounds				
2					

Comments for Sample AE05913 - Analysis \$GCMS

Westchester Dept. of Labs & Research

User: Vinci, David L

Date: 03-21-2002 Time: 15:40:34

The GCMS scan, from 35 to 550 amu, does not reveal the presence of unusual compounds. Recoveries for 27 of 43 compounds in the LCS Standard were above their acceptance ranges, while one was below. Recalculation of recovery values will be performed.

Quantitation Report (QT Reviewed)

Data File : C:\MSDCHEM\1\DATA\031702\5913.D

Acq On : 19 Mar 2002 12:18 am

Sample : 3/14/02

Misc :

MS Integration Params: BENZO.P

Quant Time: Mar 20 11:21:48 2002

Vial: 29

Operator: McLean

Inst : Instrumen

Multiplr: 1.00

Quant Results File: 5080302.RES

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

Title : Quantitation For Method 508gcscan

Last Update : Thu Mar 14 13:02:27 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	1109640	20.00	ug/L	0.00
10) Naphthalene-d8	13.03	136	3134368	20.00	ug/L	0.00
17) Acenaphthene-d10	18.13	164	1745462	20.00	ug/L	0.00
29) Phenanthrene-d10	22.49	188	2676117	20.00	ug/L	0.00
38) Chrysene-d12	30.30	240	2458182	20.00	ug/L	-0.02
44) Perylene-d12	34.18	264	1399775	20.00	ug/L	0.00

System Monitoring Compounds

37) 2,4-ddd surr	27.43	235	223924	7.01	ug/L	0.00
Spiked Amount	5.000		Recovery	=	140.20%	

Target Compounds

				Qvalue
2) N-nitroso-dimethyl amine	0.00	74	0	N.D. 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	0.00	149	0	N.D. 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	0.00	228	0	N.D. d
42) Bis (2-ethylhexyl) phthala	0.00	149	0	N.D.
43) Chrysene	0.00	228	0	N.D. 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

5913.D 5080302.M Wed Mar 20 11:23:36 2002

Quantitation Report (QT Reviewed)

Data File : C:\MSDCHEM\1\DATA\031702\5913.D Vial: 29
Acq On : 19 Mar 2002 12:18 am Operator: McLean
Sample : 3/14/02 Inst : Instrumen
Misc : Multiplr: 1.00
MS Integration Params: BENZO.P
Quant Time: Mar 20 11:21:48 2002 Quant Results File: 5080302.RES

Quant Method : C:\MSDCHEM\1\5973N\5080302.M (RTE Integrator)
Title : Quantitation For Method 508gcscan
Last Update : Thu Mar 14 13:02:27 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	0.00	252	0		N.D. 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
5913.D 5080302.M Wed Mar 20 11:23:36 2002

Data File : C:\MSDCHEM\1\DATA\031702\5913.D

Vial: 29

Acq On : 19 Mar 2002 12:18 am

Operator: McLean

Sample : 3/14/02

Inst : Instrumen

Misc :

Multiplr: 1.00

MS Integration Params: BENZO.P

Quant Time: Mar 20 11:22 2002

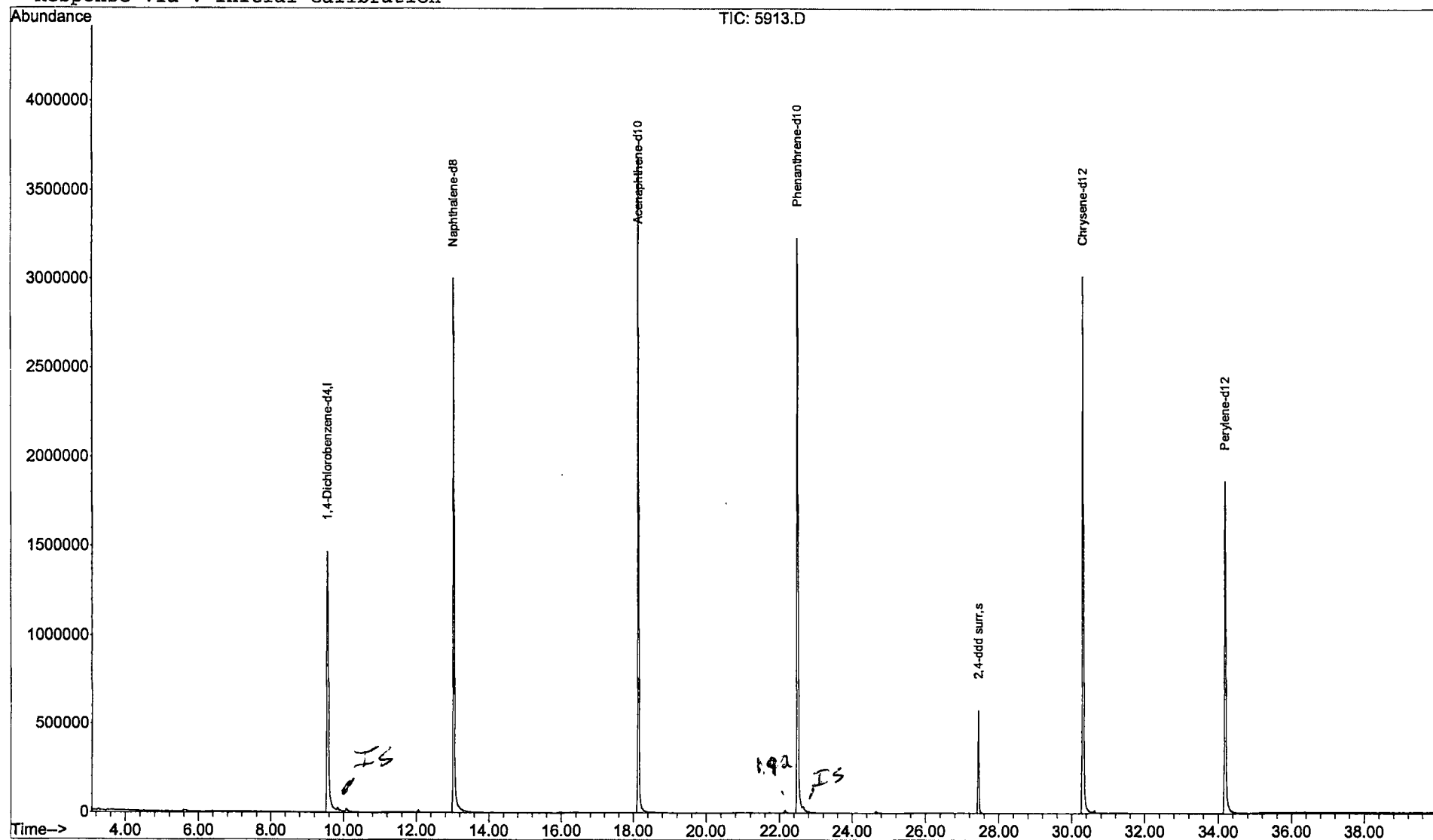
Quant Results File: 5080302.RES

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

Title : Quantitation For Method 508gcscan

Last Update : Thu Mar 14 13:02:27 2002

Response via : Initial Calibration



LSC Area Percent Report

Data File : C:\MSDCHEM\1\DATA\031702\5913.D

Acq On : 19 Mar 2002 12:18 am

Sample : 3/14/02

Misc :

MS Integration Params: LSCINT.P

Vial: 29

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

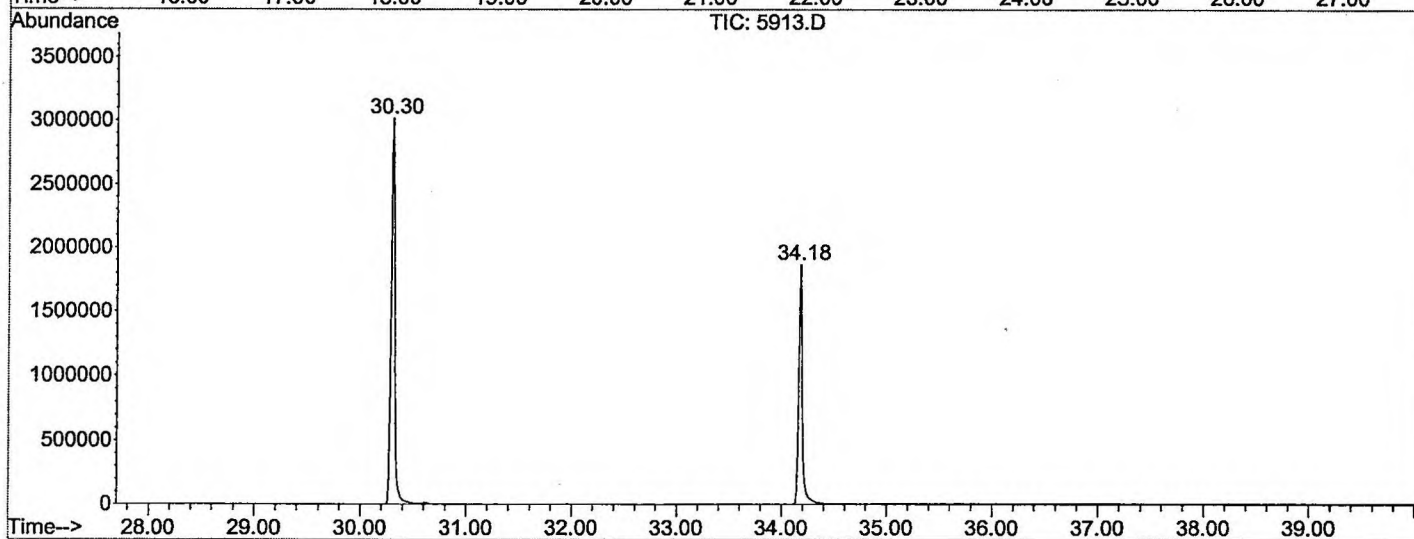
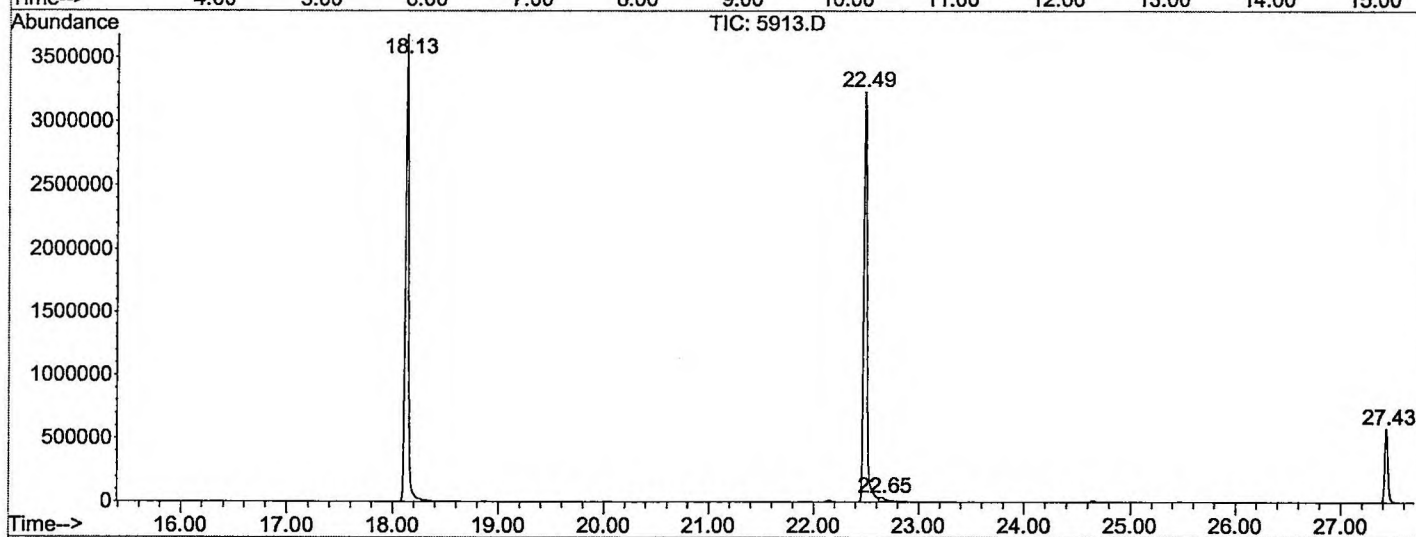
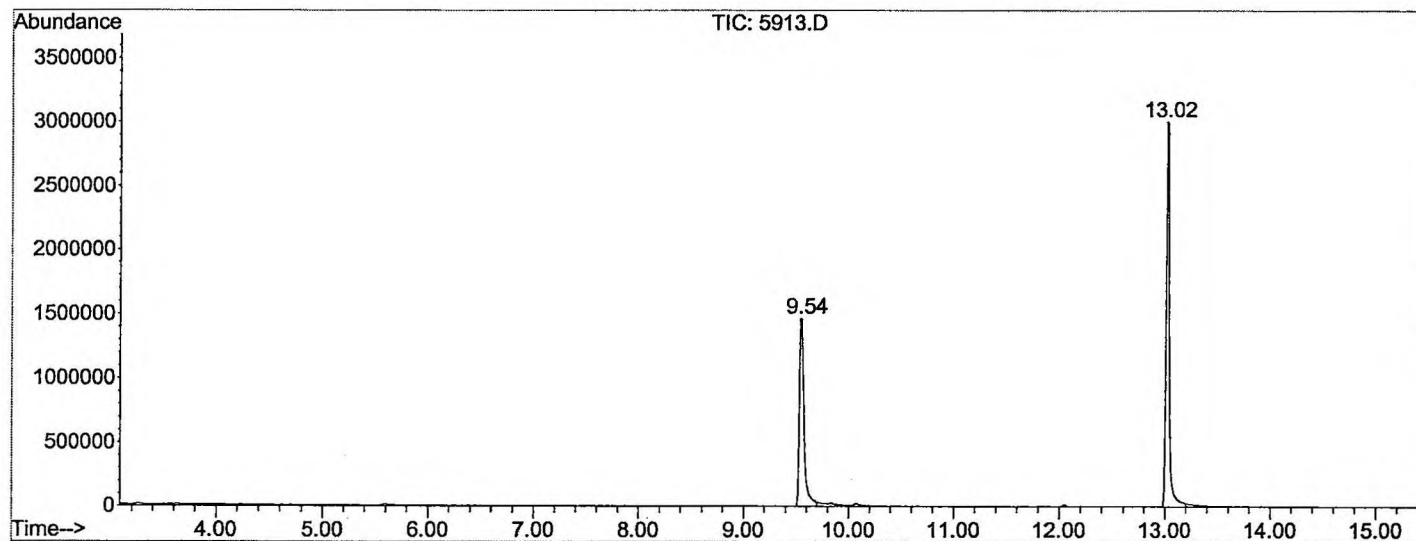
Signal : TIC

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	9.543	708	713	730	rBV	1464455	4647807	63.38%	11.927%
2	13.017	1091	1096	1116	rBV	3003177	6714430	91.56%	17.230%
3	18.133	1654	1660	1677	rBV	3681500	7253090	98.90%	18.612%
4	22.486	2135	2140	2155	rBV2	3231785	7333632	100.00%	18.818%
5	22.650	2155	2158	2172	rVB3	31689	131836	1.80%	0.338%
6	27.430	2680	2685	2694	rBV	579599	1159614	15.81%	2.976%
7	30.305	2995	3002	3025	rBV2	3016509	7159865	97.63%	18.373%
8	34.178	3422	3429	3446	rBV	1865089	4570083	62.32%	11.727%

Sum of corrected areas: 38970357

LSC Report - Integrated Chromatogram

File : C:\MSDCHEM\1\DATA\031702\5913.D
 Operator : McLean
 Acquired : 19 Mar 2002 12:18 am using AcqMethod 508SCAN
 Instrument : Instrumen
 Sample Name: 3/14/02
 Misc Info :
 Vial Number: 29
 Quant File :5080302.RES (RTE Integrator)



Tentatively Identified Compound (LSC) summary

Operator ID: McLean Date Acquired: 19 Mar 2002 12:18 am
 Data File: C:\MSDCHEM\1\DATA\031702\5913.D
 Name: 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