

User: Vinci, David L
 Westchester Dept. of Labs & Research
 Date: 04/02/2002 Time: 14:01:19

Sample: AE04382
 Analysis: \$GCMS GCMS Scan For Unknowns
 Units: ug
 Started: 4/2/02 14:00 Ended: 4/2/02 14:00 Analyst: DLV
 Page: 1

	Component Name	Viol	Result	Qualifier	MDL
1	Other unknown compounds		.		
2	Surr_2,4-DDD		119		5.0

Comments for Sample AE04382 - Analysis \$GCMS

Westchester Dept. of Labs & Research

User: Vinci, David L

Date: 04-02-2002 Time: 14:01:35

The GCMS scan, from 35 to 550 amu, does not reveal the presence of unusual compounds. The recoveries for 12 of the 43 compounds in the LCS Standard were high. New limits will be calculated.

Quantitation Report (QT Reviewed)

Data File : C:\MSDCHEM\1\DATA\032602\4382.D

Acq On : 27 Mar 2002 9:20 am

Sample : SAMPLE 4382

Misc :

MS Integration Params: BENZO.P

Quant Time: Mar 27 11:45:49 2002

Vial: 35

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 : 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	1470725	20.00	ug/L	0.00
10) Naphthalene-d8	13.03	136	4447494	20.00	ug/L	0.00
17) Acenaphthene-d10	18.14	164	2493291	20.00	ug/L	0.00
29) Phenanthrene-d10	22.50	188	3830051	20.00	ug/L	0.00
38) Chrysene-d12	30.31	240	3307310	20.00	ug/L	-0.01
44) Perylene-d12	34.18	264	1901207	20.00	ug/L	0.00

System Monitoring Compounds

37) 2,4-DDD (SURR)	27.43	235	272432	5.96	ug/L	0.00
Spiked Amount	5.000		Recovery	=	119.20%	

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	24.64	149	235611	1.16	ug/L 97
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) antracene	0.00	228	0	N.D.	d
42) Bis (2-ethylhexyl) phthala	0.00	149	0	N.D.	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

4382.D 5080302.M Wed Mar 27 11:47:40 2002

Quantitation Report

(QT Reviewed)

Data File : C:\MSDCHEM\1\DATA\032602\4382.D

Vial: 35

Acq On : 27 Mar 2002 9:20 am

Operator: McLean

Sample : SAMPLE 4382

Inst : Instrumen

Misc :

Multiplr: 1.00

MS Integration Params: BENZO.P

Quant Time: Mar 27 11:45:49 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

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

4382.D 5080302.M

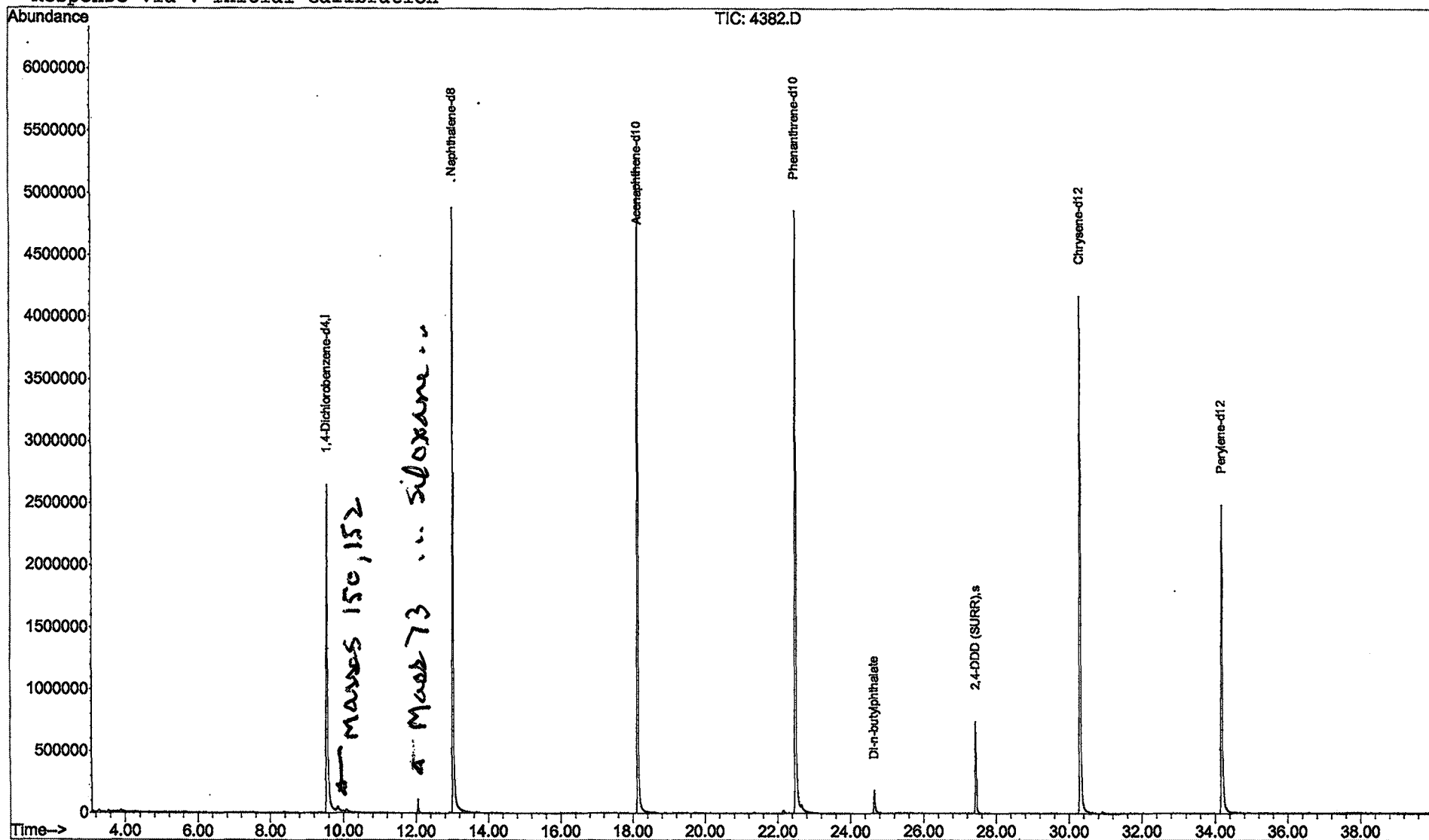
Wed Mar 27 11:47:40 2002

Data File : C:\MSDCHEM\1\DATA\032602\4382.D
 Acq On : 27 Mar 2002 9:20 am
 Sample : SAMPLE 4382
 Misc :
 MS Integration Params: BENZO.P
 Quant Time: Mar 27 11:47 2002

Vial: 35
 Operator: McLean
 Inst : Instrumen
 Multiplr: 1.00

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\032602\4382.D

Acq On : 27 Mar 2002 9:20 am

Sample : SAMPLE 4382

Misc :

MS Integration Params: LSCINT.P

Vial: 35

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

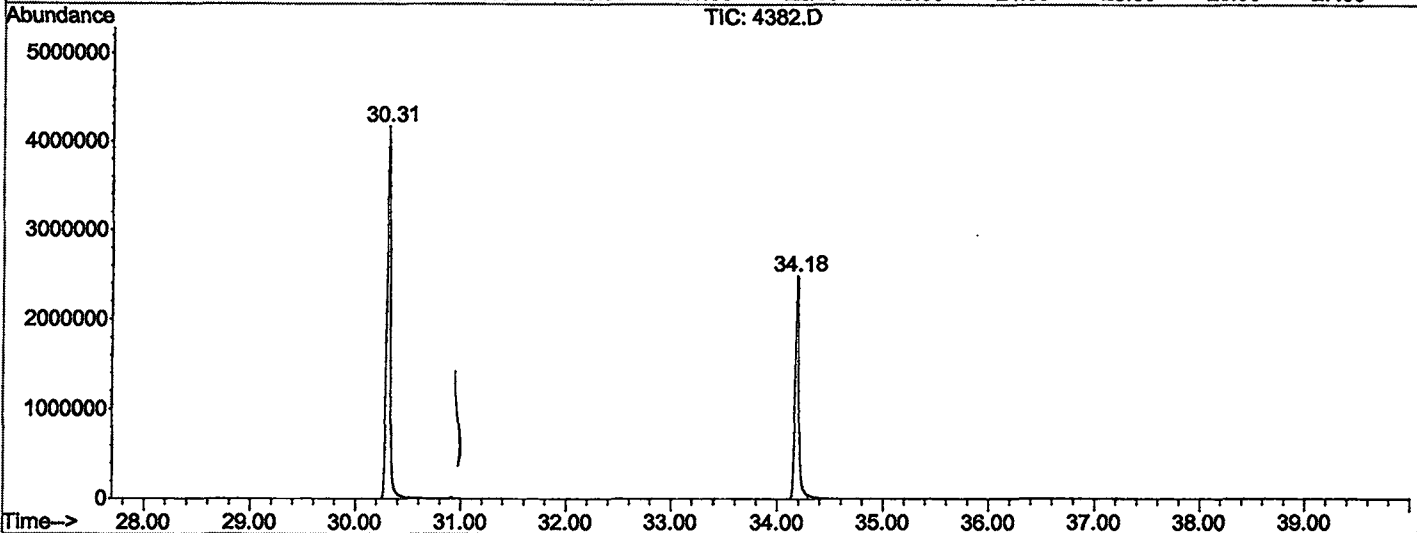
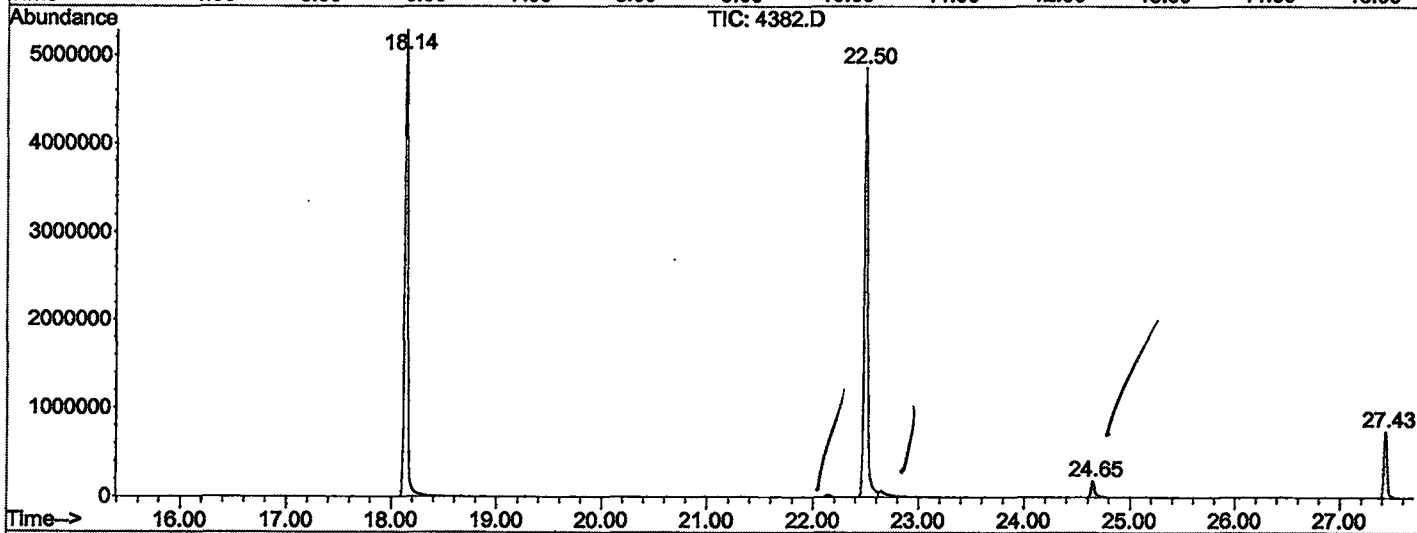
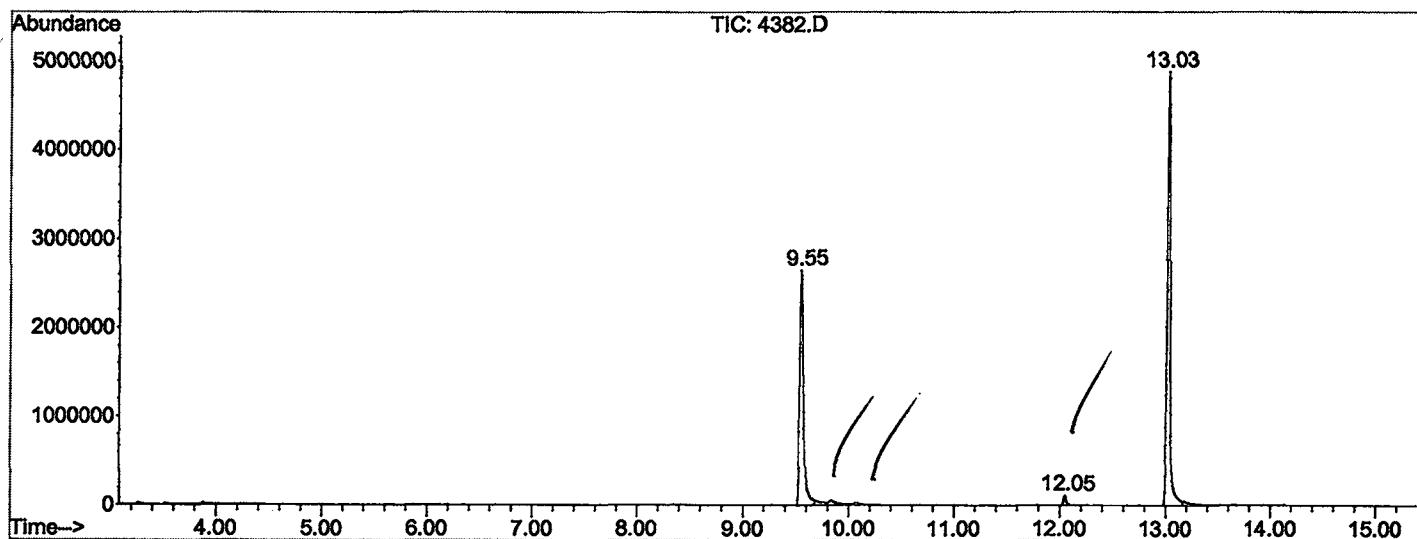
Signal : TIC

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	1123	rBV	2648851	6846642	62.94%	12.016%
2	12.051	1521	1527	1535	rBV	109973	174159	1.60%	0.306%
3	13.027	1685	1693	1716	rBV	4882622	9936843	91.35%	17.439%
4	18.138	2553	2563	2587	rBV	5278053	10810489	99.38%	18.972%
5	22.498	3295	3305	3325	rBV2	4859786	10877471	100.00%	19.090%
6	24.648	3664	3671	3682	rBV	183851	427384	3.93%	0.750%
7	27.433	4137	4145	4159	rBV2	742074	1549134	14.24%	2.719%
8	30.312	4624	4635	4664	rBV2	4173430	9979096	91.74%	17.513%
9	34.184	5283	5294	5312	rBV2	2490143	6380212	58.66%	11.197%

Sum of corrected areas: 56981430

LSC Report - Integrated Chromatogram

File : C:\MSDCHEM\1\DATA\032602\4382.D
 Operator : McLean
 Acquired : 27 Mar 2002 9:20 am using AcqMethod 508SCAN
 Instrument : Instrumen
 Sample Name: SAMPLE 4382
 Misc Info :
 Vial Number: 35
 Quant File :5080302.RES (RTE Integrator)



Tentatively Identified Compound (LSC) summary

Operator ID: McLean Date Acquired: 27 Mar 2002 9:20 am
 Data File: C:\MSDCHEM\1\DATA\032602\4382.D
 Name: SAMPLE 4382
 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