

User: Vinci, David L
 Westchester Dept. of Labs & Research
 Date: 04/11/2002 Time: 16:11:02

Sample: AE06851
 Analysis: \$GCMS GCMS Scan For Unknowns
 Units: ug
 Started: 4/11/02 16:10 Ended: 4/11/02 16:10 Analyst: DLV
 Page: 1

	Component Name	Vol	Result	Qualifier	MDL
1	Other unknown compounds		.		
2	Other unknown compounds		0.000000		0.0

Comments for Sample AE06851 - Analysis \$GCMS

Westchester Dept. of Labs & Research

User: Vinci, David L

Date: 04-11-2002 Time: 16:11:19

The GCMS scan, from 35 to 550 amu, does not reveal the presence of unusual compounds. The recoveries for 8 of the 43 compounds in the LCS Standard were high.

Quantitation Report

(QT Reviewed)

Data File : C:\MSDCHEM\1\DATA\040502\6851.D

Acq On : 7 Apr 2002 7:06 am

Sample : SAMPLE 6851 3/22/02

Misc :

MS Integration Params: BENZO.P

Quant Time: Apr 10 07:43:02 2002

Vial: 44

Operator: Bohlk

Inst : Instrumen

Multiplr: 1.00

Quant Results File: 5080402.RES

Quant Method : C:\MSDCHEM\1\5973N\5080402.M (RTE Integrator)

Title : Quantitation For Method 508gcscan

Last Update : Mon Apr 08 06:47:22 2002

Response via : Initial Calibration

DataAcq Meth : 508SCAN

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	9.46	152	995848	20.00	ug/L	-0.01
10) Naphthalene-d8	12.93	136	4487106	20.00	ug/L	-0.01
17) Acenaphthene-d10	18.05	164	2626597	20.00	ug/L	0.00
29) Phenanthrene-d10	22.40	188	3768684	20.00	ug/L	-0.02
38) Chrysene-d12	30.22	240	2916441	20.00	ug/L	-0.02
44) Perylene-d12	34.11	264	1798118	20.00	ug/L	-0.01

System Monitoring Compounds

37) 2,4-DDD (SURR)	27.36	235	276732	7.40	ug/L	0.00
Spiked Amount	5.000		Recovery	= 148.00%		

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	30.85	149	9645	0.10 ug/L	84
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

6851.D 5080402.M Wed Apr 10 07:44:58 2002

Quantitation Report

(QT Reviewed)

Data File : C:\MSDCHEM\1\DATA\040502\6851.D

Vial: 44

Acq On : 7 Apr 2002 7:06 am

Operator: Bohlk

Sample : SAMPLE 6851 3/22/02

Inst : Instrumen

Misc :

Multiplr: 1.00

MS Integration Params: BENZO.P

Quant Time: Apr 10 07:43:02 2002

Quant Results File: 5080402.RES

Quant Method : C:\MSDCHEM\1\5973N\5080402.M (RTE Integrator)

Title : Quantitation For Method 508gcscan

Last Update : Mon Apr 08 06:47:22 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

6851.D 5080402.M Wed Apr 10 07:44:58 2002

Page 2

Data File : C:\MSDCHEM\1\DATA\040502\6851.D

Vial: 44

Acq On : 7 Apr 2002 7:06 am

Operator: Bohlk

Sample : SAMPLE 6851 3/22/02

Inst : Instrumen

Misc :

Multiplr: 1.00

MS Integration Params: BENZO.P

Quant Time: Apr 10 10:44 2002

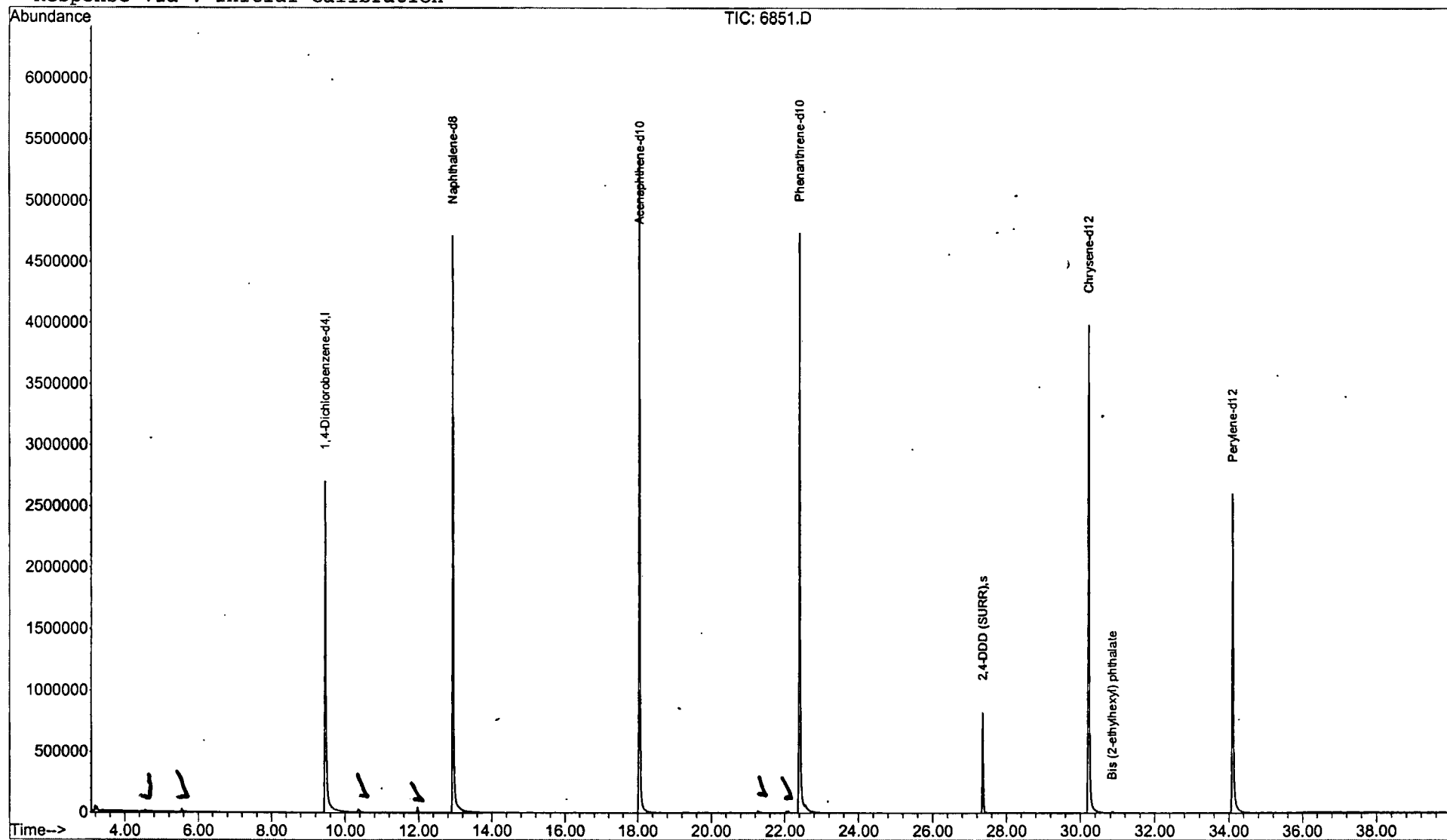
Quant Results File: 5080402.RES

Method : C:\MSDCHEM\1\5973N\5080402.M (RTE Integrator)

Title : Quantitation For Method 508gcscan

Last Update : Mon Apr 08 06:47:22 2002

Response via : Initial Calibration



LSC Area Percent Report

Data File : C:\MSDCHEM\1\DATA\040502\6851.D
Acq On : 7 Apr 2002 7:06 am
Sample : SAMPLE 6851 3/22/02
Misc :
MS Integration Params: LSCINT.P

Vial: 44
Operator: Bohlk
Inst : Instrumen
Multiplr: 1.00

Method : C:\MSDCHEM\1\5973N\5080402.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

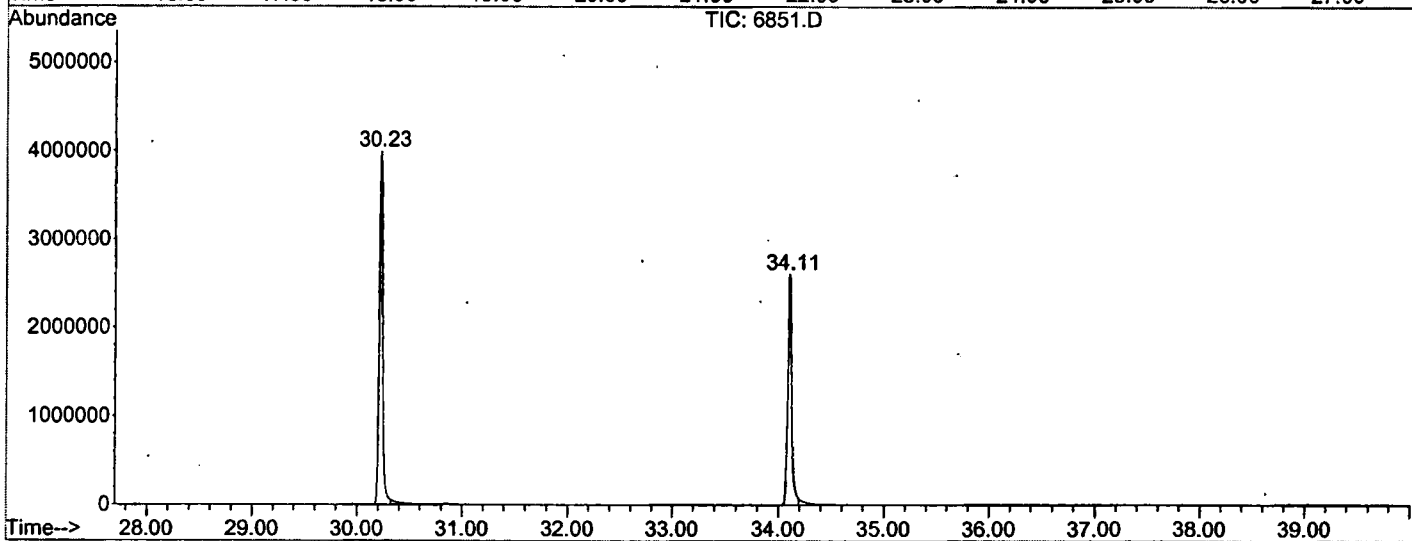
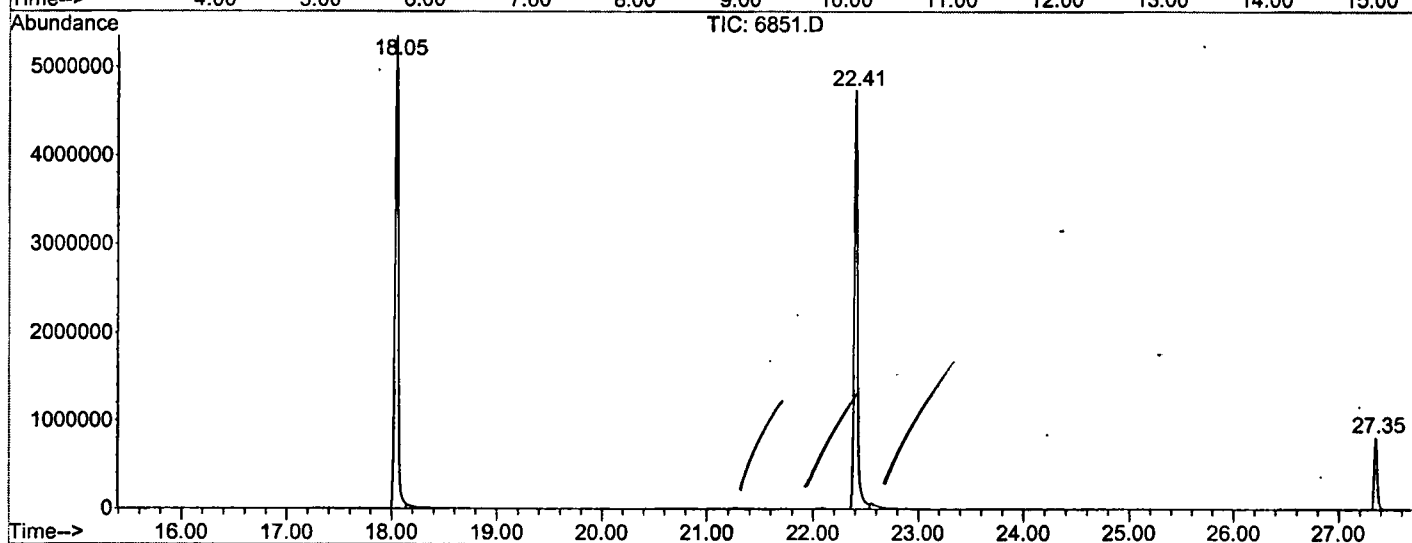
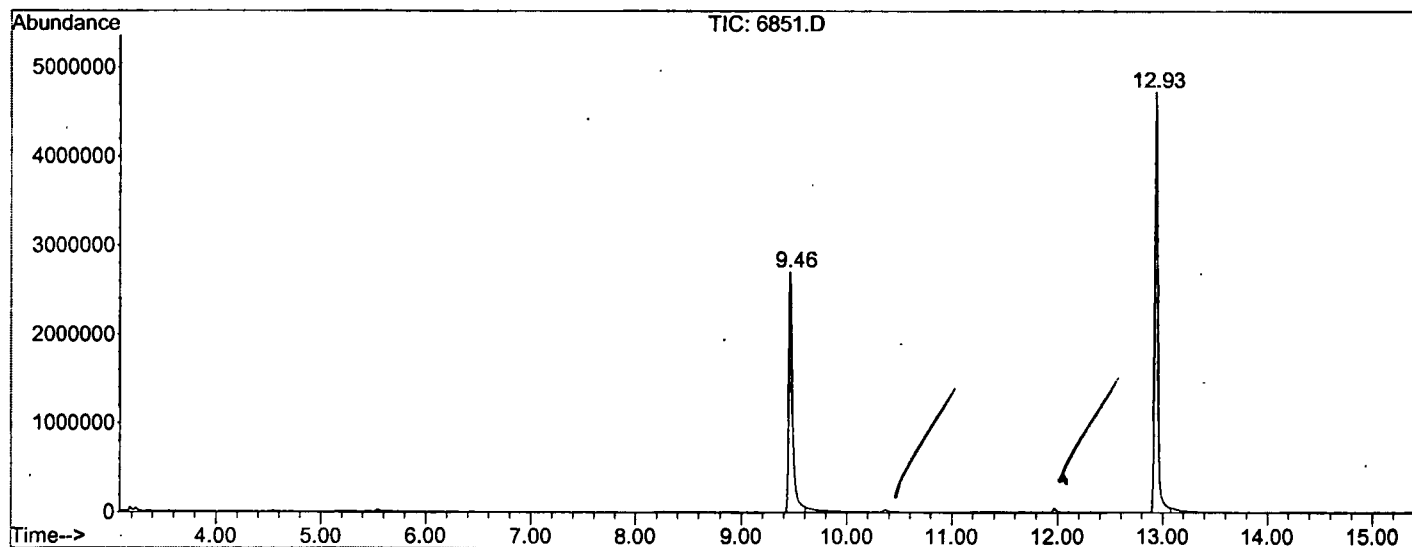
Signal : TIC

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	9.460	1079	1086	1110	rBV	2699830	6707865	59.51%	11.846%
2	12.933	1669	1677	1701	rBV	4714751	9936344	88.15%	17.547%
3	18.050	2538	2548	2563	rBV	5349053	11271809	100.00%	19.906%
4	22.410	3280	3290	3312	rBV2	4739509	11021412	97.78%	19.463%
5	27.351	4124	4131	4144	rBV	816880	1611526	14.30%	2.846%
6	30.230	4610	4621	4637	rBV2	3989949	9391255	83.32%	16.585%
7	34.108	5269	5281	5296	rBV2	2602512	6686168	59.32%	11.808%

Sum of corrected areas: 56626379

LSC Report - Integrated Chromatogram

File : C:\MSDCHEM\1\DATA\040502\6851.D
Operator : Bohlk
Acquired : 7 Apr 2002 7:06 am using AcqMethod 508SCAN
Instrument : Instrumen
Sample Name: SAMPLE 6851 3/22/02
Misc Info :
Vial Number: 44
Quant File :5080402.RES (RTE Integrator)



Tentatively Identified Compound (LSC) summary

Operator ID: Bohlk Date Acquired: 7 Apr 2002 7:06 am
 Data File: C:\MSDCHEM\1\DATA\040502\6851.D
 Name: SAMPLE 6851 3/22/02
 Misc:
 Method: C:\MSDCHEM\1\5973N\5080402.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
