

Jrd

User: Vinci, David L
 Westchester Dept. of Labs & Research
 Date: 03/22/2002 Time: 18:19:51

Sample: AE03078
 Analysis: \$GCMS GCMS Scan For Unknowns
 Units: ug
 Started: 3/22/02 18:19 Ended: 3/22/02 18:19 Analyst: DLV
 Page: 1

	Component Name	Viol	Result	Qualifier	MDL
1	Other unknown compounds				
2					

Comments for Sample AE03078 - Analysis \$GCMS

Westchester Dept. of Labs & Research

User: Vinci, David L

Date: 03-22-2002 Time: 18:19:54

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

Quantitation Report (QT Reviewed)

Data File : C:\MSDCHEM\1\DATA\031802\3078.D

Vial: 9

Acq On : 19 Mar 2002 8:45 am

Operator: McLean

Sample : GC/MS SCAN 3/6 FB

Inst : Instrumen

Misc :

Multiplr: 1.00

MS Integration Params: BENZO.P

Quant Time: Mar 19 21:34:20 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

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	9.55	150	1280057	20.00	ug/L	0.00
10) Naphthalene-d8	13.03	136	3538043	20.00	ug/L	0.00
17) Acenaphthene-d10	18.13	164	1972638	20.00	ug/L	0.00
29) Phenanthrene-d10	22.50	188	3013416	20.00	ug/L	0.00
38) Chrysene-d12	30.30	240	2790862	20.00	ug/L	-0.02
44) Perylene-d12	34.18	264	1451998	20.00	ug/L	0.00

System Monitoring Compounds

37) 2,4-ddd surr	27.43	235	249254	6.93	ug/L	0.00
Spiked Amount	5.000		Recovery	=	138.60%	

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	18.13	165	251870	11.01 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	24.65	149	41551	0.26 ug/L	98
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	30.91	149	19678	0.18 ug/L	89
43) Chrysene	30.31	228	6930	0.05 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

3078.D 5080302.M Tue Mar 19 21:34:43 2002

Quantitation Report

(QT Reviewed)

Data File : C:\MSDCHEM\1\DATA\031802\3078.D

Acq On : 19 Mar 2002 8:45 am

Sample : GC/MS SCAN 3/6 FB

Misc :

MS Integration Params: BENZO.P

Quant Time: Mar 19 21:34:20 2002

Vial: 9

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

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

3078.D 5080302.M Tue Mar 19 21:34:43 2002

Data File : C:\MSDCHEM\1\DATA\031802\3078.D

Acq On : 19 Mar 2002 8:45 am

Sample : GC/MS SCAN 3/6 FB

Misc :

MS Integration Params: BENZO.P

Quant Time: Mar 19 21:34 2002

Vial: 9

Operator: McLean

Inst : Instrumen

Multiplr: 1.00

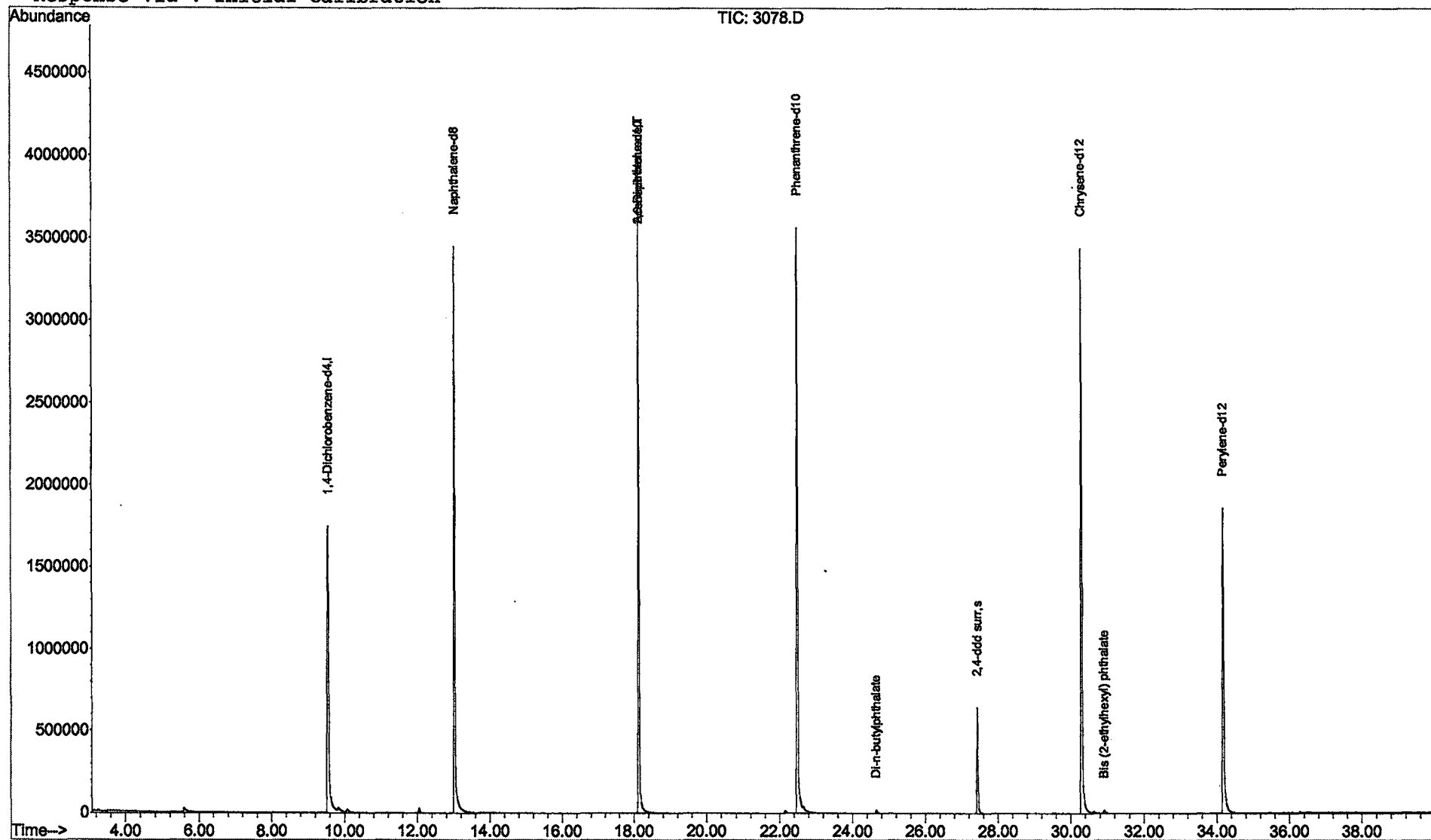
Quant Results File: 5080302.RES

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

Title : Quantitation For Method 508gcscan

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

Response via : Initial Calibration



LSC Area Percent Report

Data File : C:\MSDCHEM\1\DATA\031802\3078.D
Acq On : 19 Mar 2002 8:45 am
Sample : GC/MS SCAN 3/6 FB
Misc :
MS Integration Params: LSCINT.P

Vial: 9
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

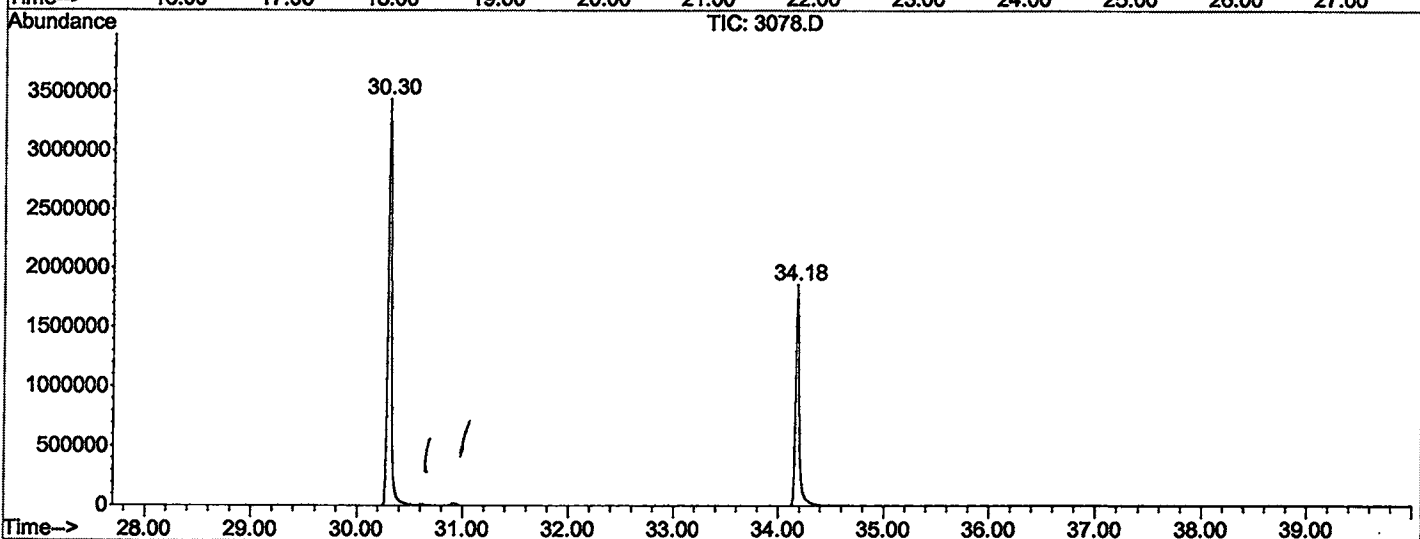
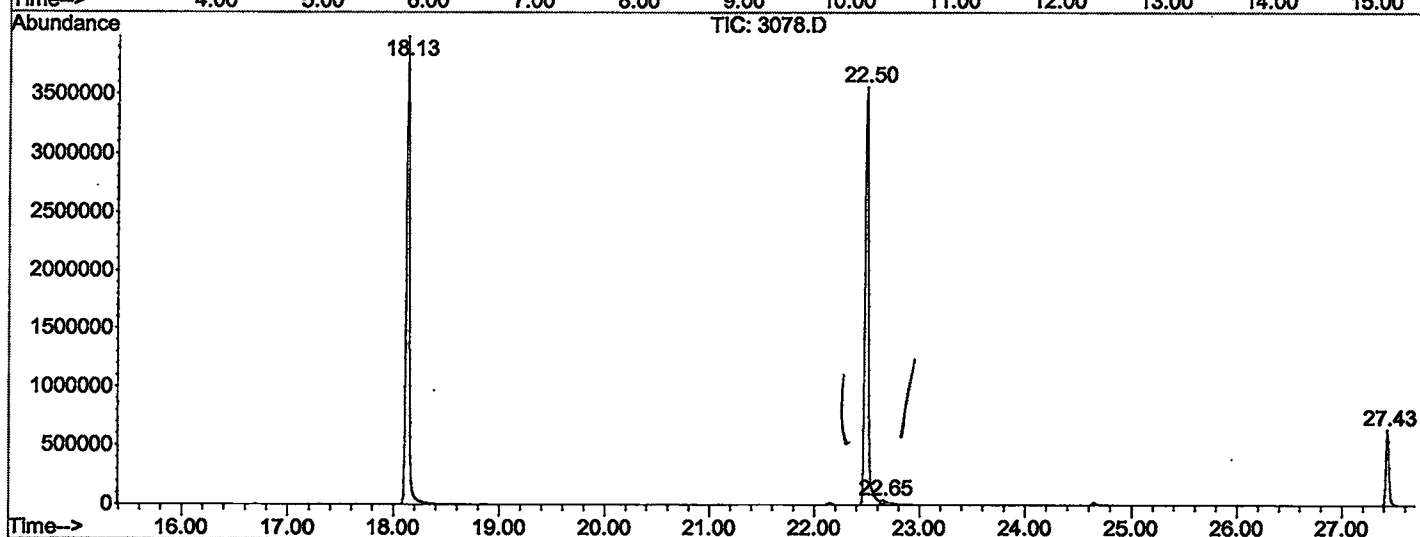
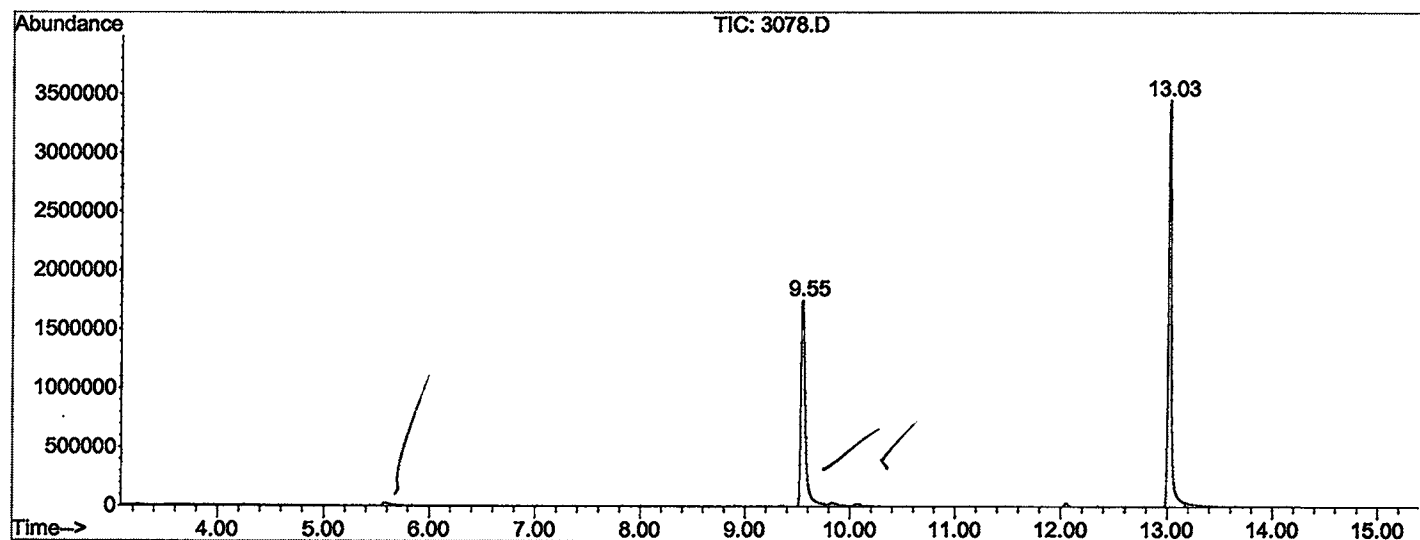
Signal : TIC

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	9.552	709	714	738	rBV	1743629	5324780	64.73%	12.327%
2	13.026	1091	1097	1113	rBV	3447662	7423439	90.24%	17.186%
3	18.133	1654	1660	1679	rBV	3986468	8097923	98.44%	18.747%
4	22.495	2135	2141	2155	rBV2	3561890	8226575	100.00%	19.045%
5	22.650	2155	2158	2165	rVB3	27472	87277	1.06%	0.202%
6	27.430	2681	2685	2699	rBV	642690	1270084	15.44%	2.940%
7	30.305	2996	3002	3023	rBV2	3438499	8059163	97.96%	18.657%
8	34.178	3423	3429	3446	rBV2	1859651	4706151	57.21%	10.895%

Sum of corrected areas: 43195392

LSC Report - Integrated Chromatogram

File : C:\MSDCHEM\1\DATA\031802\3078.D
 Operator : McLean
 Acquired : 19 Mar 2002 8:45 am using AcqMethod 508SCAN
 Instrument : Instrumen
 Sample Name: GC/MS SCAN 3/6 FB
 Misc Info :
 Vial Number: 9
 Quant File :5080302.RES (RTE Integrator)



Tentatively Identified Compound (LSC) summary

Operator ID: McLean Date Acquired: 19 Mar 2002 8:45 am
 Data File: C:\MSDCHEM\1\DATA\031802\3078.D
 Name: GC/MS SCAN 3/6 FB
 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
