

DLV

User: Vinci, David L
 Westchester Dept. of Labs & Research
 Date: 03/22/2002 Time: 18:30:17

Sample: AE03183
 Analysis: \$GCMS GCMS Scan For Unknowns
 Units: ug
 Started: 3/22/02 18:29 Ended: 3/22/02 18:29 Analyst: DLV
 Page: 1

	Component Name	Viol	Result	Qualifier	MDL
1	Other unknown compounds				
2					

Comments for Sample AE03183 - Analysis \$GCMS

Westchester Dept. of Labs & Research

User: Vinci, David L

Date: 03-22-2002 Time: 18:30:21

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\3183.D

Vial: 15

Acq On : 19 Mar 2002 1:38 pm

Operator: McLean

Sample : GC/MS SCAN 3/6

Inst : Instrumen

Misc :

Multiplr: 1.00

MS Integration Params: BENZO.P

Quant Time: Mar 19 21:39:45 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	9.54	150	1127625	20.00	ug/L	0.00
10) Naphthalene-d8	13.02	136	3145417	20.00	ug/L	0.00
17) Acenaphthene-d10	18.13	164	1756229	20.00	ug/L	0.00
29) Phenanthrene-d10	22.49	188	2671808	20.00	ug/L	0.00
38) Chrysene-d12	30.30	240	2484195	20.00	ug/L	-0.02
44) Perylene-d12	34.18	264	1312685	20.00	ug/L	0.00

System Monitoring Compounds

37) 2,4-ddd surr	27.43	235	238646	7.49	ug/L	0.00
Spiked Amount	5.000		Recovery	=	149.80%	

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	18.13	165	223648	10.98	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	42784	0.30	ug/L	95
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	30.30	228	5707	0.05	ug/L	81
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

3183.D 5080302.M Tue Mar 19 21:40:26 2002

Quantitation Report

(QT Reviewed)

Data File : C:\MSDCHEM\1\DATA\031802\3183.D

Vial: 15

Acq On : 19 Mar 2002 1:38 pm

Operator: McLean

Sample : GC/MS SCAN 3/6

Inst : Instrumen

Misc :

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MS Integration Params: BENZO.P

Quant Time: Mar 19 21:39:45 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

3183.D 5080302.M Tue Mar 19 21:40:27 2002

Page 2

Data File : C:\MSDCHEM\1\DATA\031802\3183.D

Vial: 15

Acq On : 19 Mar 2002 1:38 pm

Operator: McLean

Sample : GC/MS SCAN 3/6

Inst : Instrumen

Misc :

Multiplr: 1.00

MS Integration Params: BENZO.P

Quant Time: Mar 19 21:40 2002

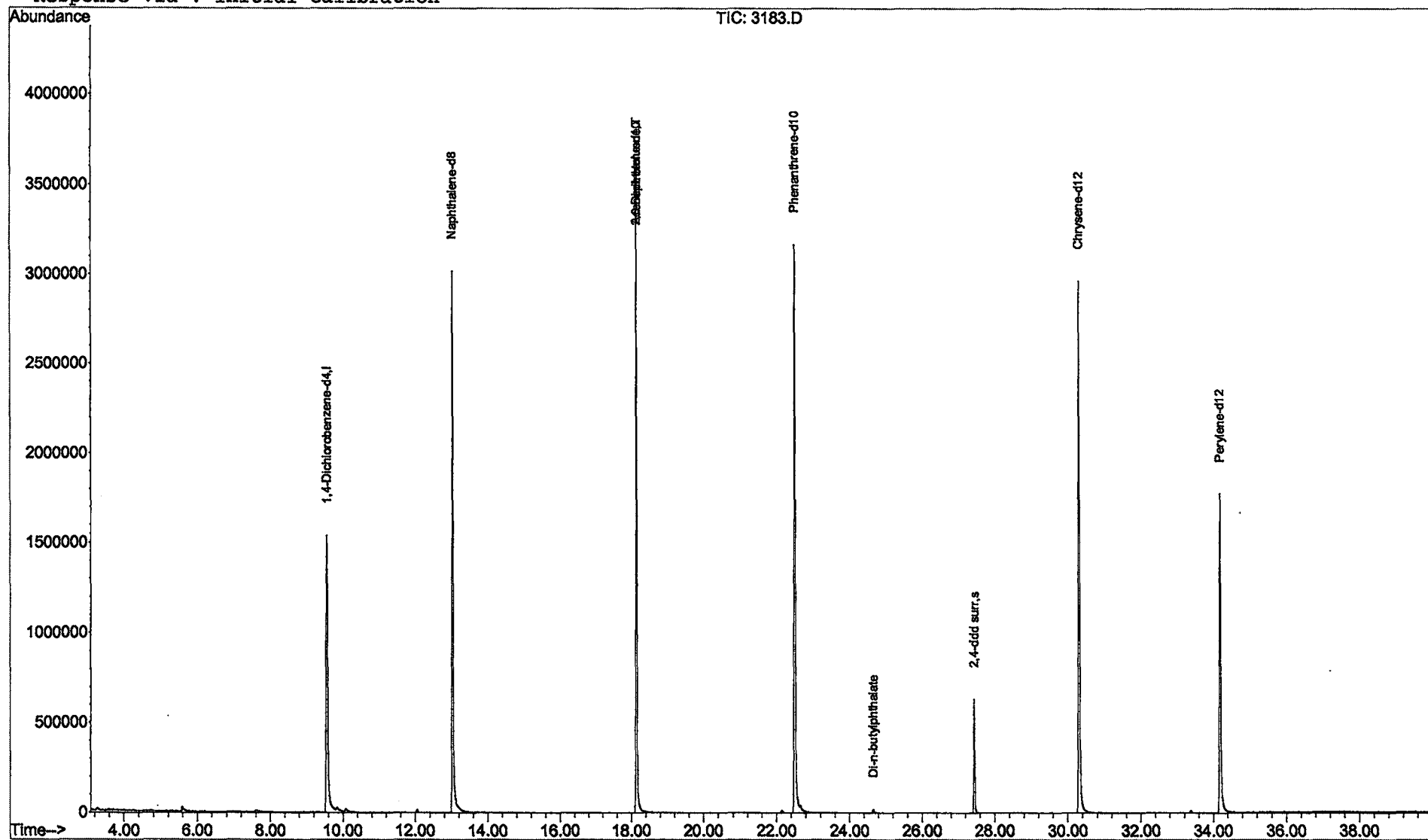
Quant Results File: 5080302.RES

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\031802\3183.D

Acq On : 19 Mar 2002 1:38 pm

Sample : GC/MS SCAN 3/6

Misc :

MS Integration Params: LSCINT.P

Vial: 15

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

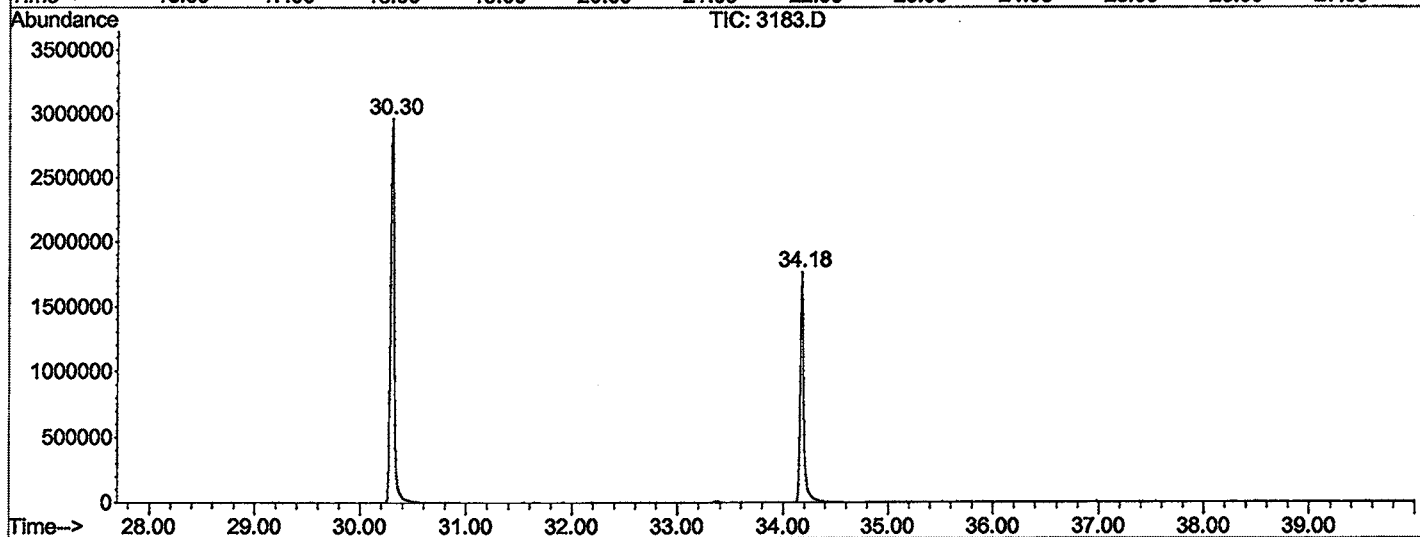
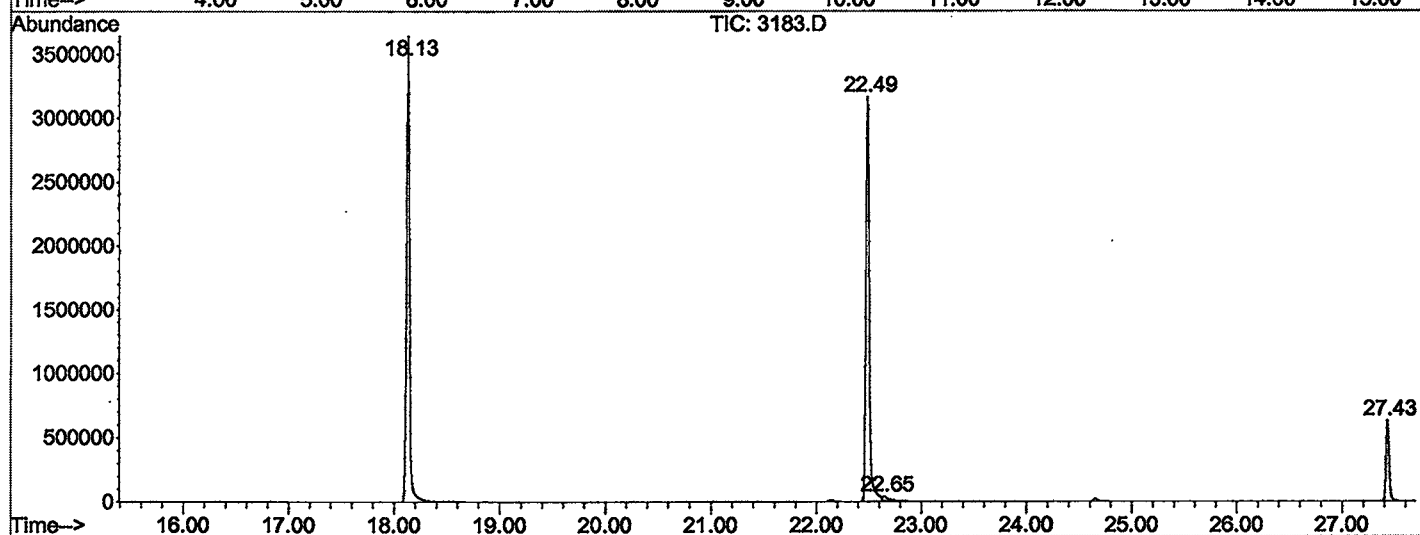
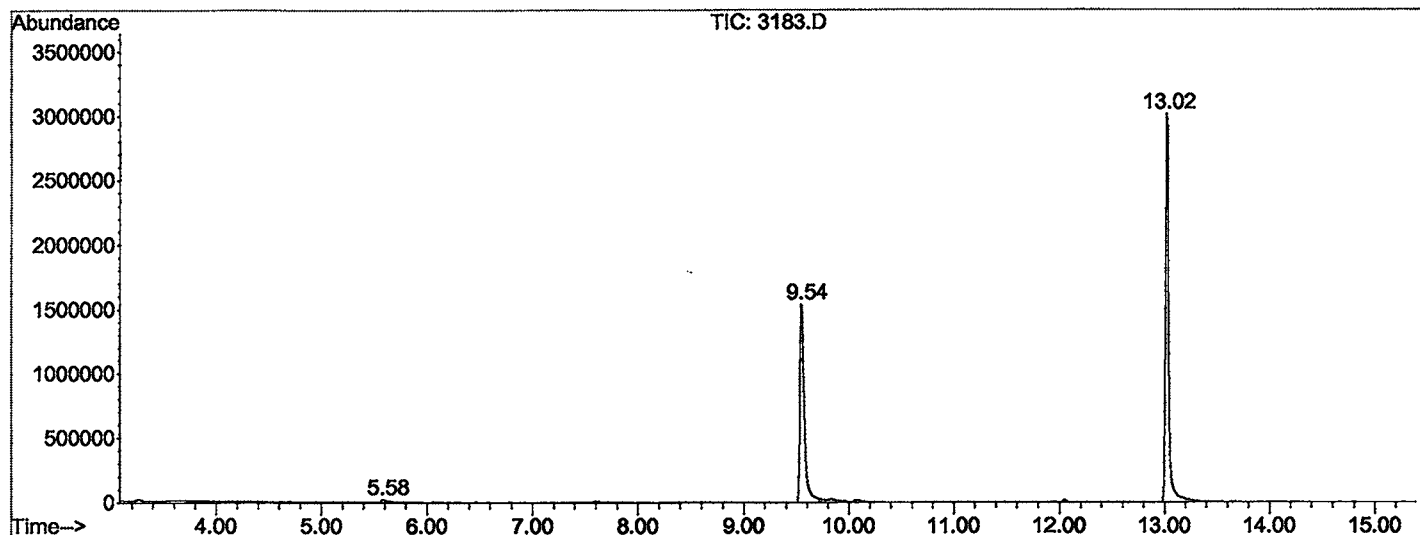
Signal : TIC

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	5.579	273	276	286	rBV2	22817	79663	1.09%	0.204%
2	9.543	709	713	732	rBV	1539817	4664697	63.63%	11.961%
3	13.017	1091	1096	1129	rBV	3017066	6824463	93.09%	17.498%
4	18.133	1654	1660	1677	rBV	3640654	7229514	98.62%	18.537%
5	22.486	2135	2140	2155	rBV2	3166087	7330653	100.00%	18.796%
6	22.650	2156	2158	2170	rVB2	36176	122126	1.67%	0.313%
7	27.430	2680	2685	2697	rBV	632100	1248973	17.04%	3.202%
8	30.305	2996	3002	3026	rBV2	2960658	7142882	97.44%	18.315%
9	34.178	3422	3429	3452	rBV	1775565	4357835	59.45%	11.174%

Sum of corrected areas: 39000806

LSC Report - Integrated Chromatogram

File : C:\MSDCHEM\1\DATA\031802\3183.D
 Operator : McLean
 Acquired : 19 Mar 2002 1:38 pm using AcqMethod 508SCAN
 Instrument : Instrumen
 Sample Name: GC/MS SCAN 3/6
 Misc Info :
 Vial Number: 15
 Quant File :5080302.RES (RTE Integrator)



Tentatively Identified Compound (LSC) summary

• Operator ID: McLean Date Acquired: 19 Mar 2002 1:38 pm
 Data File: C:\MSDCHEM\1\DATA\031802\3183.D
 Name: GC/MS SCAN 3/6
 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