User: Vinci, David L
Westchester Dept. of Labs & Research
Date: 04/29/2002 Time: 08:50:09

Sample: AE08936
Analysis: $GCMS GCMS Scan For Unknowns
Units: ug
Started: 4/29/02 08:49 Ended: 4/29/02 08:49 Analyst: DLV
Page: 1

	Component Name	Viol	Result	Qualifier	MDL
1	Other unknown compounds		.		
2	Surr_2,4-DDD		70		10.0

Comments for Sample AE08936 - Analysis $GCMS

Westchester Dept. of Labs & Research
User: Vinci, David L
Date: 04-29-2002 Time: 08:51:30

The GCMS scan, from 35 to 550 amu, does not reveal the presence of unusual compounds.

Quantitation Report (QT Reviewed)

Data File : C:\MSDCHEM\1\DATA\042202\8936.D Vial: 9
Acq On : 22 Apr 2002 5:43 pm Operator: Bohlk
Sample : SAMPLE 8936 4/15/02 Inst : Instrumen
Misc : Multiplr: 1.00
MS Integration Params: BENZO.P
Quant Time: Apr 25 14:31:22 2002 Quant Results File: 5080402BBC.RES

Quant Method : C:\MSDCHEM\1\5973N\5080402BBC.M (RTE Integrator)
Title : Quantitation For Method 508gcscan
Last Update : Wed Apr 24 11:07:03 2002
Response via : Initial Calibration
DataAcq Meth : 508SCAN

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	9.41	152	350145	40.00	ug/L	0.00
10) Naphthalene-d8	12.88	136	1666663	40.00	ug/L	0.00
17) Acenaphthene-d10	17.99	164	943737	40.00	ug/L	0.00
30) Phenanthrene-d10	22.34	188	1317501	40.00	ug/L	0.00
39) Chrysene-d12	30.14	240	971849	40.00	ug/L	-0.01
45) Perylene-d12	34.01	264	533227	40.00	ug/L	0.00
System Monitoring Compounds						
28) TCMX (SURR)	19.95	244	26252	6.97	ug/L	0.00
Spiked Amount 10.000			Recovery = 69.70%			
38) 2,4-DDD (SURR)	0.00	235	0	0.00	ug/L	
Spiked Amount 10.000			Recovery = 0.00%			
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.		
29) Azobenzen/ 1,2-Diphenylhyd	0.00	77	0	N.D.		
31) N-Nitrosodiphenylamine	0.00	169	0	N.D.		
32) 4-Bromophenyl-phenyl ether	0.00	248	0	N.D.		
33) Hexachlorobenzene	0.00	284	0	N.D.		
34) Phenanthrene	0.00	178	0	N.D.		
35) Anthracene	0.00	178	0	N.D.		
36) Di-n-butylphthalate	0.00	149	0	N.D.		
37) Fluoranthene	0.00	202	0	N.D.		
40) Pyrene	0.00	202	0	N.D.		
41) Butylbenzylphthalate	0.00	149	0	N.D.		
42) Benzo (a) antracene	0.00	228	0	N.D.		
43) Bis (2-ethylhexyl) phthala	0.00	149	0	N.D.		
44) Chrysene	0.00	228	0	N.D.		

(#) = qualifier out of range (m) = manual integration

Quantitation Report (QT Reviewed)

Data File : C:\MSDCHEM\1\DATA\042202\8936.D Vial: 9
Acq On : 22 Apr 2002 5:43 pm Operator: Bohlk
Sample : SAMPLE 8936 4/15/02 Inst : Instrumen
Misc : Multiplr: 1.00
MS Integration Params: BENZO.P
Quant Time: Apr 25 14:31:22 2002 Quant Results File: 5080402BBC.RES

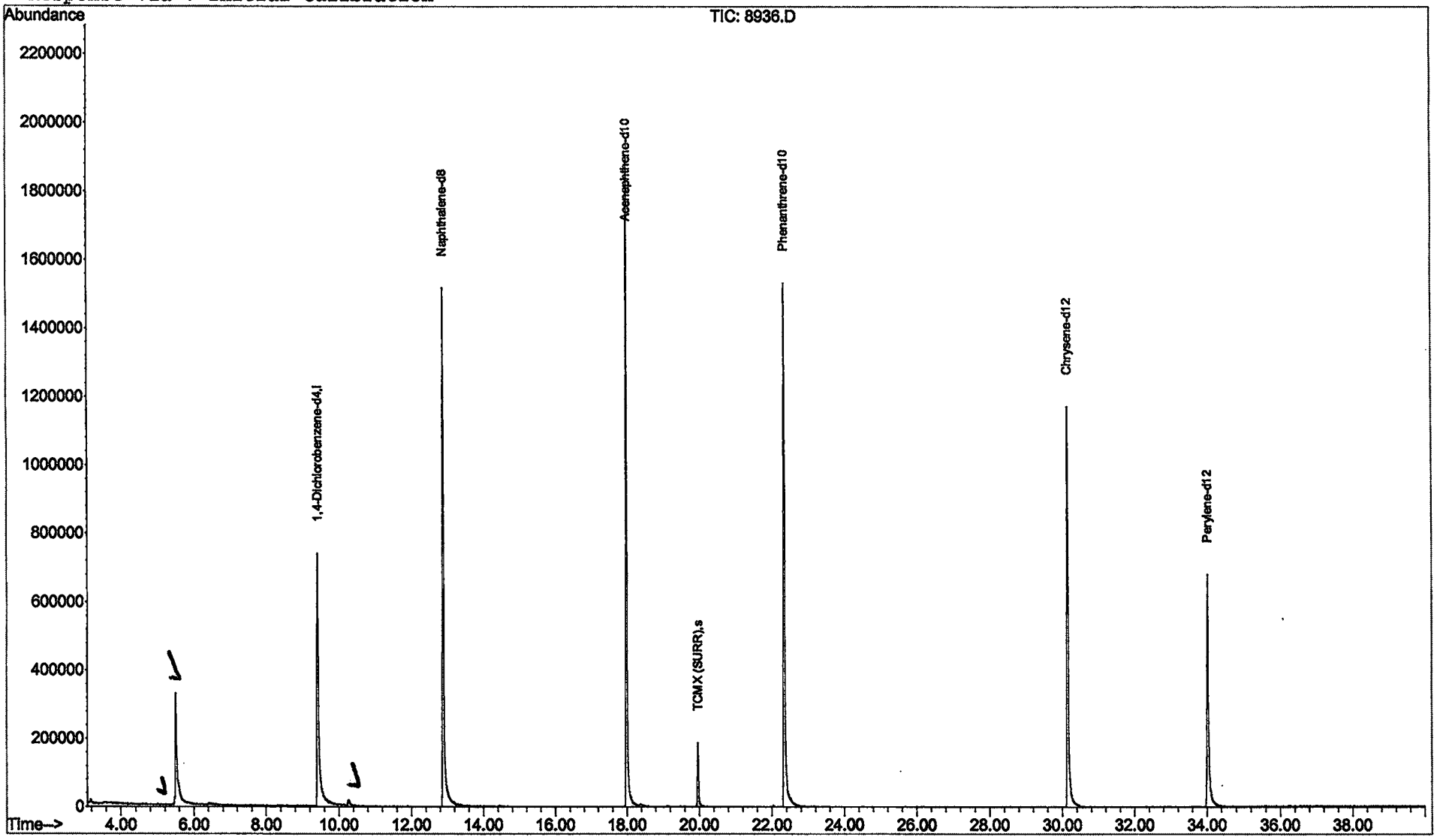
Quant Method : C:\MSDCHEM\1\5973N\5080402BBC.M (RTE Integrator)
Title : Quantitation For Method 508gcscan
Last Update : Wed Apr 24 11:07:03 2002
Response via : Initial Calibration
DataAcq Meth : 508SCAN

Compound	R.T.	QIon	Response	Conc Unit	Qvalue
46) Di-n-Octylphthalate	0.00	149	0	N.D.	
47) Benzo (b) fluoranthene	0.00	252	0	N.D.	
48) Benzo (k) fluoranthene	0.00	252	0	N.D.	
49) Benzo (a) pyrene	0.00	252	0	N.D.	
50) Indeno (1,2,3-cd) pyrene	0.00	276	0	N.D.	
51) Dibenz (a,h) anthracene	0.00	278	0	N.D.	
52) Benzo (g,h,i) perylene	0.00	276	0	N.D.	

(#) = qualifier out of range (m) = manual integration
8936.D 5080402BBC.M Thu Apr 25 14:32:03 2002 Page 2

Data File : C:\MSDCHEM\1\DATA\042202\8936.D Vial: 9
Acq On : 22 Apr 2002 5:43 pm Operator: Bohlk
Sample : SAMPLE 8936 4/15/02 Inst : Instrumen
Misc : Multiplr: 1.00
MS Integration Params: BENZO.P
Quant Time: Apr 25 14:31 2002 Quant Results File: 5080402BBC.RES

Method : C:\MSDCHEM\1\5973N\5080402BBC.M (RTE Integrator)
Title : Quantitation For Method 508gcscan
Last Update : Wed Apr 24 11:07:03 2002
Response via : Initial Calibration



LSC Area Percent Report

Data File : C:\MSDCHEM\1\DATA\042202\8936.D Vial: 9
Acq On : 22 Apr 2002 5:43 pm Operator: Bohlk
Sample : SAMPLE 8936 4/15/02 Inst : Instrumen
Misc : Multiplr: 1.00
MS Integration Params: LSCINT.P

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

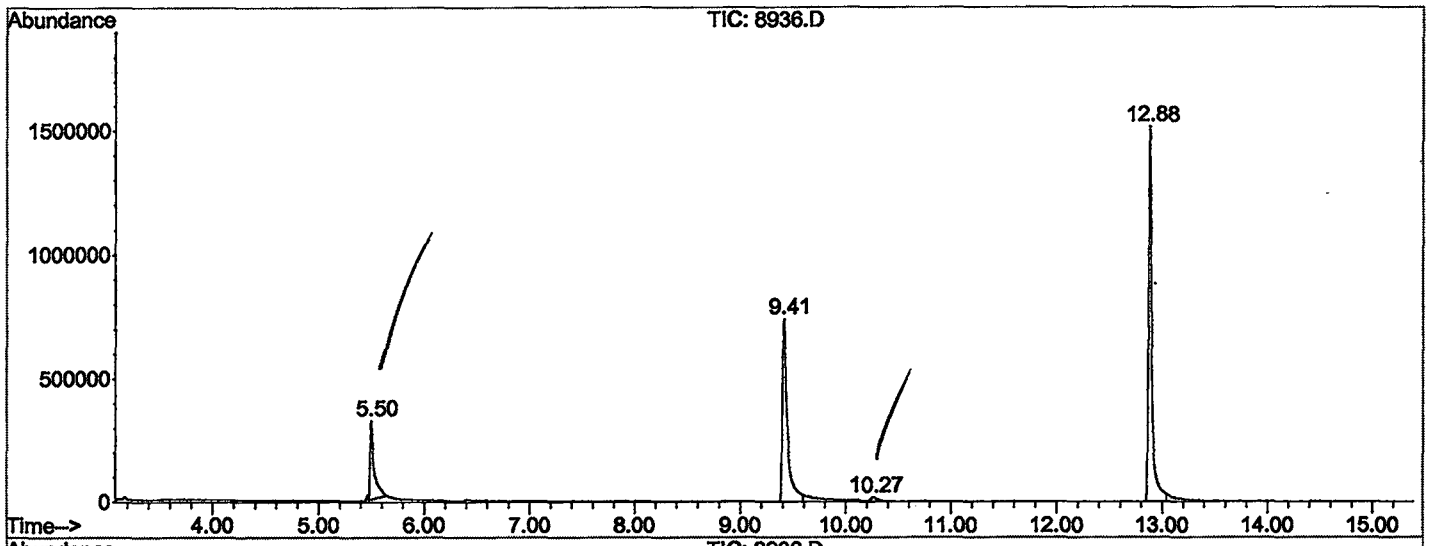
Signal : TIC

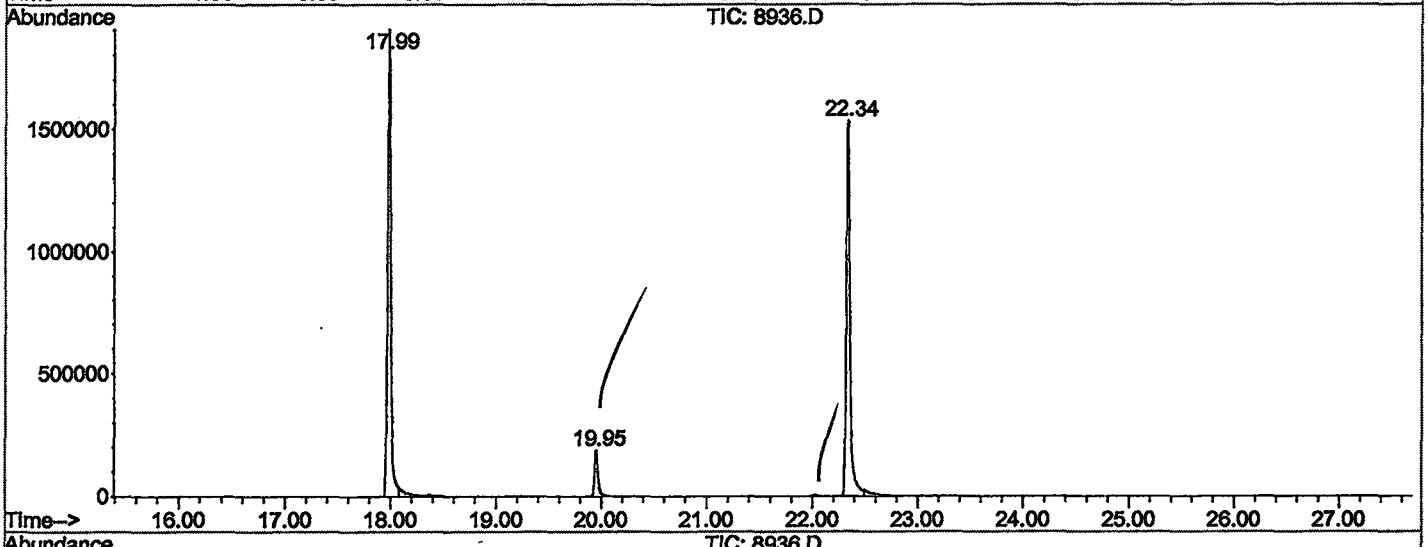
peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	5.500	408	412	436	rVB	318229	843086	21.21%	4.227%
2	9.413	1072	1078	1108	rBV	739751	2379382	59.85%	11.929%
3	10.265	1216	1223	1229	rBV4	16320	49154	1.24%	0.246%
4	12.880	1661	1668	1695	rBV	1516341	3571866	89.85%	17.908%
5	17.986	2529	2537	2553	rBV	1903739	3975554	100.00%	19.931%
6	19.954	2864	2872	2891	rBV3	187424	411845	10.36%	2.065%
7	22.339	3270	3278	3304	rBV2	1533386	3783984	95.18%	18.971%
8	30.142	4597	4606	4631	rBV2	1171721	3005173	75.59%	15.066%
9	34.014	5256	5265	5286	rBV	680717	1926091	48.45%	9.656%

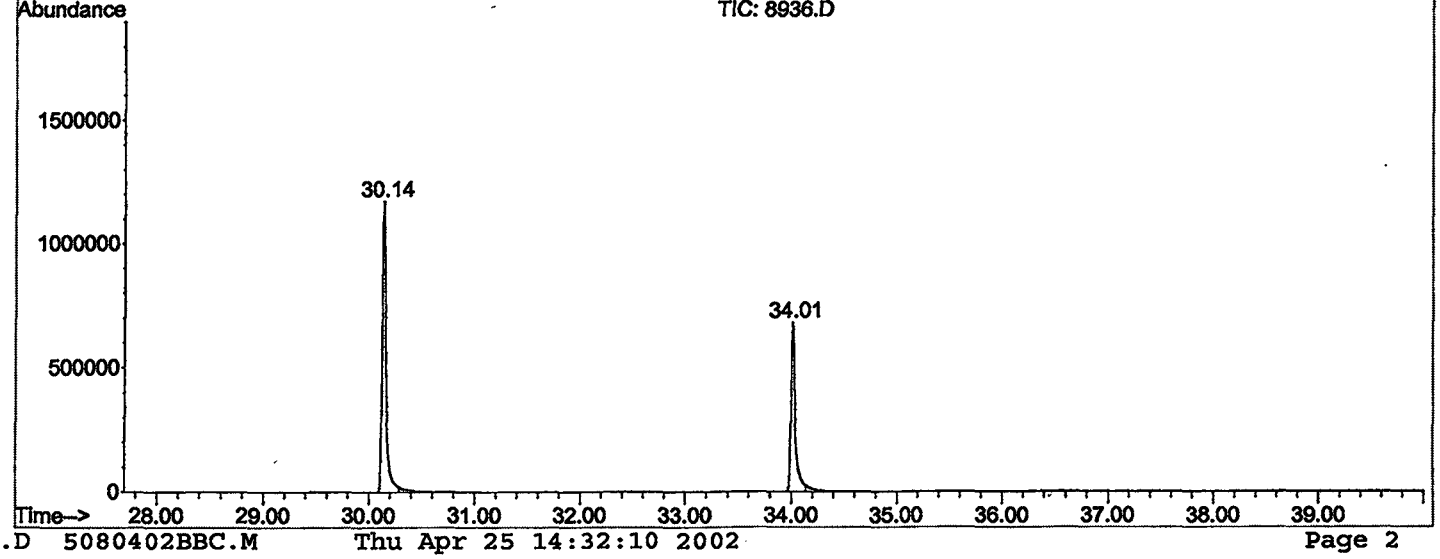
Sum of corrected areas: 19946135

LSC Report - Integrated Chromatogram

File : C:\MSDCHEM\1\DATA\042202\8936.D
Operator : Bohlk
Acquired : 22 Apr 2002 5:43 pm using AcqMethod 508SCAN
Instrument : Instrumen
Sample Name: SAMPLE 8936 4/15/02
Misc Info :
Vial Number: 9
Quant File :5080402BBC.RES (RTE Integrator)







Library Search Compound Report

Data File : C:\MSDCHEM\1\DATA\042202\8936.D Vial: 9
Acq On : 22 Apr 2002 5:43 pm Operator: Bohlk
Sample : SAMPLE 8936 4/15/02 Inst : Instrumen
Misc : Multiplr: 1.00
MS Integration Params: LSCINT.P

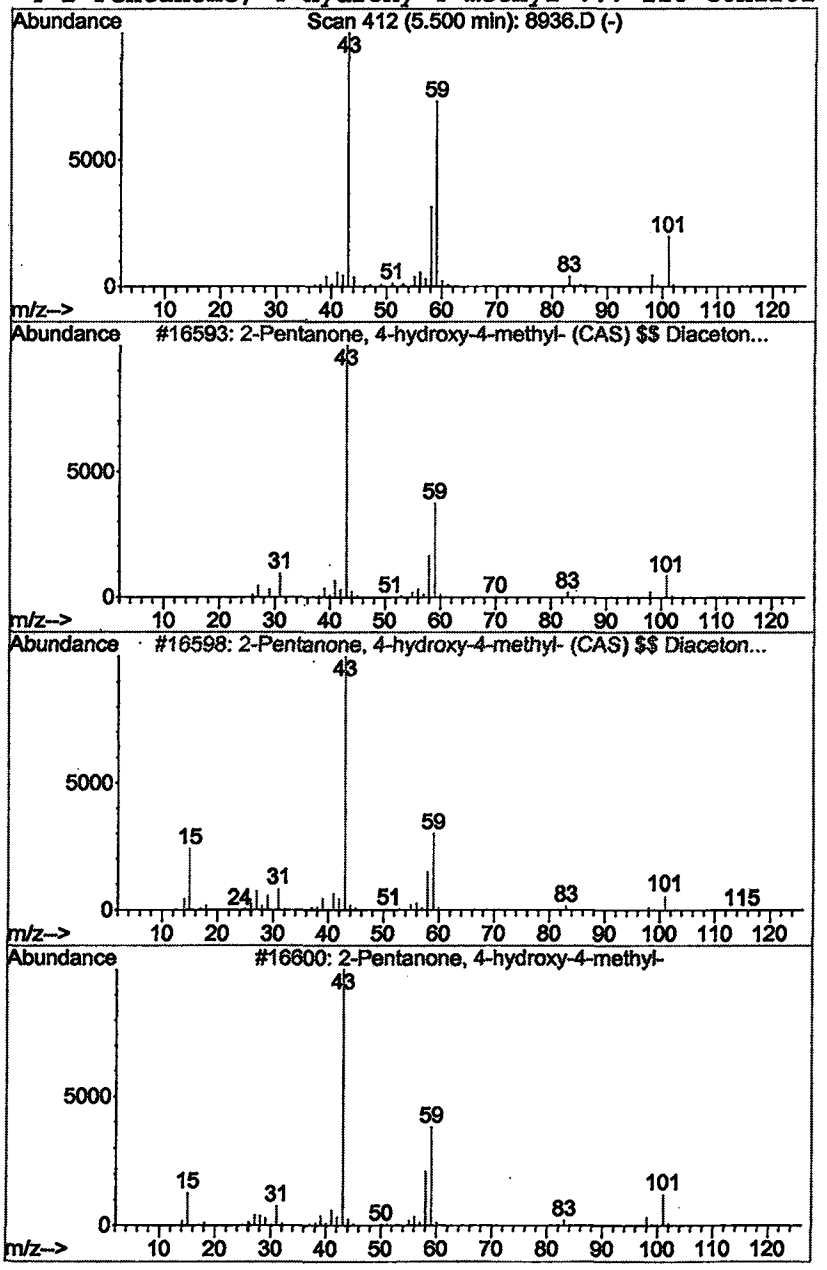
Quant Method : C:\MSDCHEM\1\5973N\5080402BBC.M (RTE Integrator)
Title : Quantitation For Method 508gcscan
Library : C:\DATABASE\WILEY7N.L

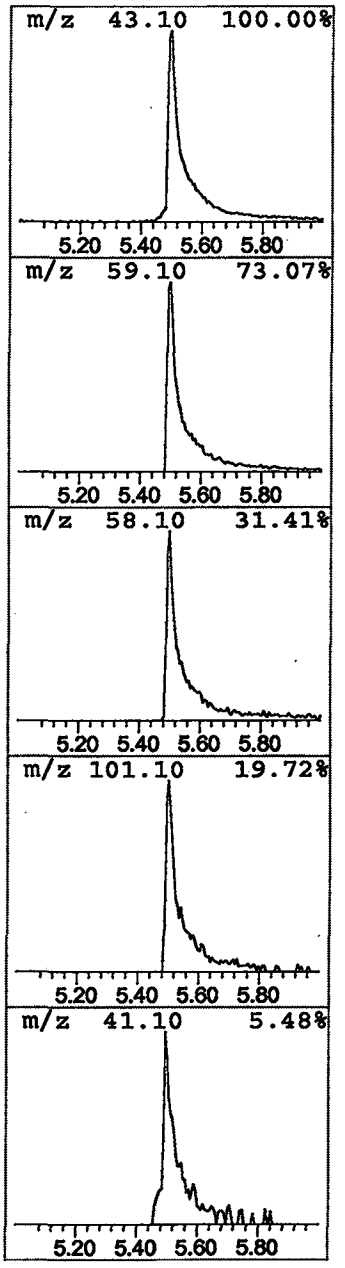
****

Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.50	14.17 ug/L	843086	1,4-Dichlorobenzene-d4	9.41

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Pentanone, 4-hydroxy-4-methyl-...	116	C6H12O2	000123-42-2	83
2	2-Pentanone, 4-hydroxy-4-methyl-...	116	C6H12O2	000123-42-2	83
3	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	83
4	2-Pentanone, 4-hydroxy-4-methyl-...	116	C6H12O2	000123-42-2	72





Library Search Compound Report

Data File : C:\MSDCHEM\1\DATA\042202\8936.D Vial: 9
Acq On : 22 Apr 2002 5:43 pm Operator: Bohlk
Sample : SAMPLE 8936 4/15/02 Inst : Instrumen
Misc : Multiplr: 1.00
MS Integration Params: LSCINT.P

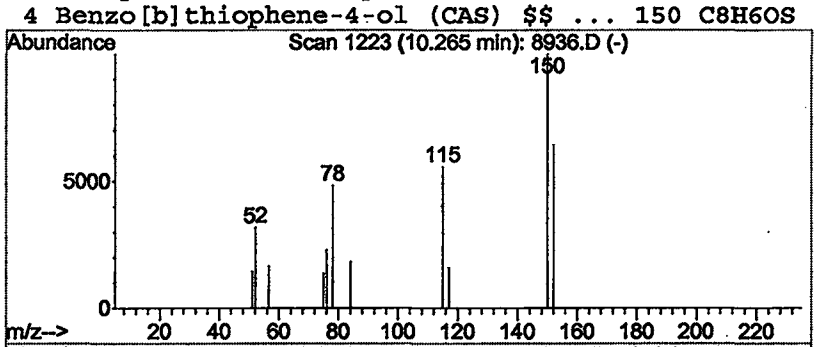
Quant Method : C:\MSDCHEM\1\5973N\5080402BBC.M (RTE Integrator)
Title : Quantitation For Method 508gcscan
Library : C:\DATABASE\WILEY7N.L

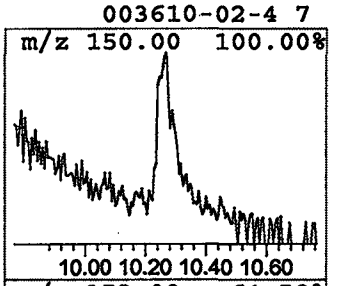
****

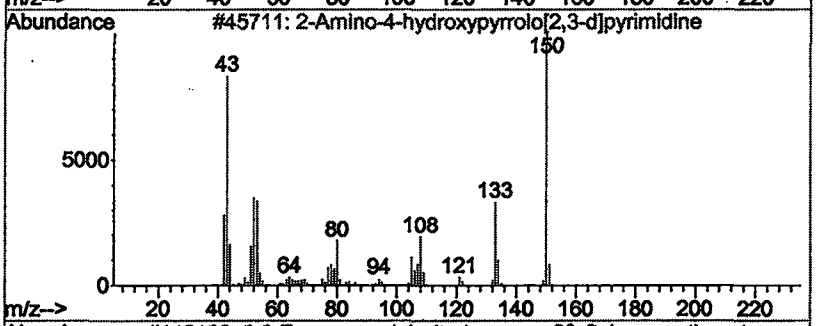
Peak Number 2 2-Amino-4-hydroxypyrrolo[2,... Concentration Rank 2

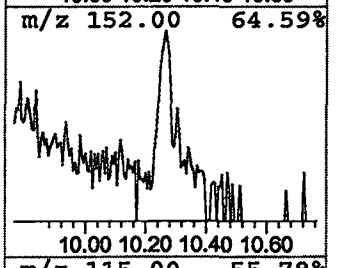
R.T.	EstConc	Area	Relative to ISTD	R.T.
10.27	0.83 ug/L	49154	1,4-Dichlorobenzene-d4	9.41

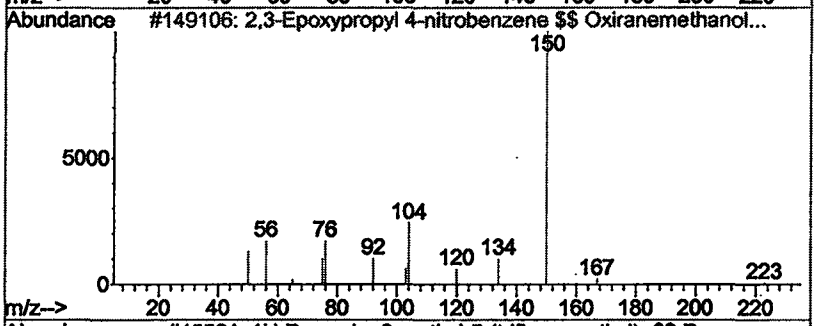
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Amino-4-hydroxypyrrolo[2,3-d]p...	150	C6H6N4O	000000-00-0	9
2	2,3-Epoxypropyl 4-nitrobenzene $...	223	C10H9NO5	013318-11-1	9
3	1H-Pyrazole, 3-methyl-5-(trifluo...	150	C5H5F3N2	010010-93-2	7
4	Benzo[b]thiophene-4-ol (CAS) $$ ...	150	C8H6OS	003610-02-4	7

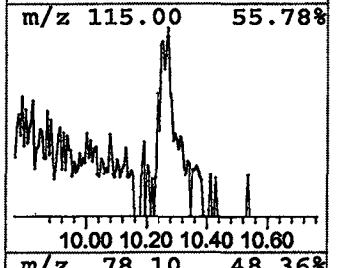


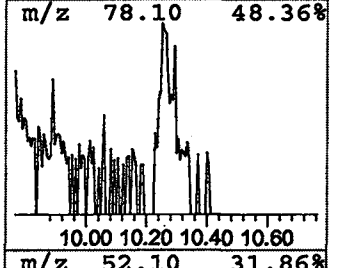


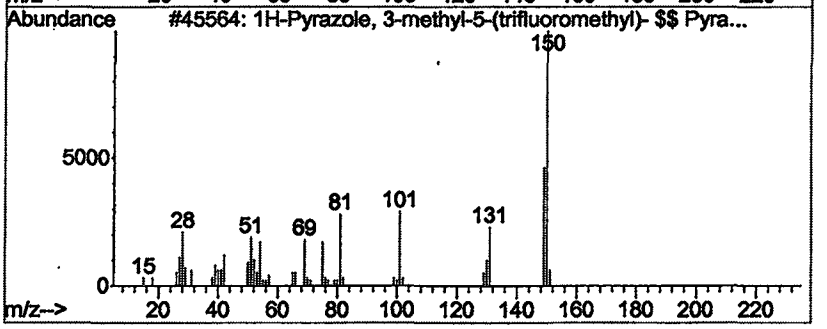


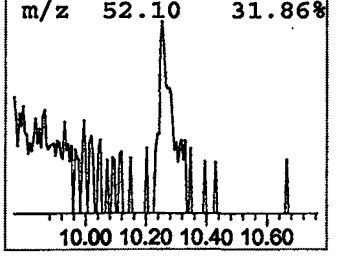












Tentatively Identified Compound (LSC) summary

Operator ID: Bohlk Date Acquired: 22 Apr 2002 5:43 pm
Data File: C:\MSDCHEM\1\DATA\042202\8936.D
Name: SAMPLE 8936 4/15/02
Misc:
Method: C:\MSDCHEM\1\5973N\5080402BBC.M (RTE Integrator)
Title: Quantitation For Method 508gcscan
Library Searched: C:\DATABASE\WILEY7N.L

					--Internal Standard---			
TIC Top Hit name	RT	EstConc	Units	Response	#	RT	Resp	Conc
2-Pentanone, 4-hy...	5.50	14.2	ug/L	843086	1	9.41	2379380	40.0
2-Amino-4-hydroxy...	10.27	0.8	ug/L	49154	1	9.41	2379380	40.0