

User: Vinci, David L
Westchester Dept. of Labs & Research
Date: 02/11/2002 Time: 15:42:54

Sample: AD31346
Analysis: \$GCMS GCMS Scan For Unknowns
Units: ug
Started: 2/11/02 15:42 Ended: 2/11/02 15:42 Analyst: DLV
Page: 1

	Component Name	Viol	Result	Qualifier	MDL
1	Other unknown compounds		.		
2	Surr_2,4-DDD		87		5.0

Comments for Sample AD31346 - Analysis \$GCMS

Westchester Dept. of Labs & Research

User: Vinci, David L

Date: 02-11-2002 Time: 15:42:57

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

Data File : C:\HPCHEM\1\DATA\JAN0902\31346.D

Acq On : 11 Jan 2002 2:18 am

Sample : gcms scan 12/24

Misc :

MS Integration Params: GOOFY.P

Quant Time: Feb 5 16:40 2002

Vial: 12

Operator: 209 GALOS

Inst : GC/MS 209

Multiplr: 1.00

Quant Results File: GCMS0103.RES

Quant Method : C:\HPCHEM\1\METHODS\GCMS0103.M (RTE Integrator)

Title : 209 625/TCLP calibration table

Last Update : Thu Dec 13 14:40:46 2001

Response via : Initial Calibration

DataAcq Meth : GCMS0103

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	11.53	152	624987	20.00	ug/l	0.00
10) Naphthalene-d8	15.12	136	2455525	20.00	ug/l	0.00
17) Acenaphthene-d10	20.39	164	1250682	20.00	ug/l	0.00
29) Phenanthrene-d10	24.80	188	1751643	20.00	ug/l	0.00
38) Chrysene-d12	32.84	240	1051570	20.00	ug/l	0.00
44) Perylene-d12	36.90	264	616028	20.00	ug/l	0.00

System Monitoring Compounds

37) 2,4-DDD	29.89	235	88146	4.38	ug/mL	-0.18
Spiked Amount	5.000		Recovery	=	87.60%	

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-Dichlorobenzene	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-Chloropropan	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)methane	0.00	93	0	N.D.
14) 1,2,4-Trichlorobenzene	0.00	180	0	N.D.
15) Naphthalene	15.18	128	909	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	20.75	165	100	N.D.
25) Diethylphthalate	21.93	149	267	N.D.
26) 4-Chlorophenyl-phenylether	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.

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

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Quantitation Report

(QT Reviewed)

Data File : C:\HPCHEM\1\DATA\JAN0902\31346.D

Vial: 12

Acq On : 11 Jan 2002 2:18 am

Operator: 209 GALOS

Sample : gcms scan 12/24

Inst : GC/MS 209

Misc :

Multiplr: 1.00

MS Integration Params: GOOFY.P

Quant Time: Feb 5 16:40 2002

Quant Results File: GCMS0103.RES

Quant Method : C:\HPCHEM\1\METHODS\GCMS0103.M (RTE Integrator)

Title : 209 625/TCLP calibration table

Last Update : Thu Dec 13 14:40:46 2001

Response via : Initial Calibration

DataAcq Meth : GCMS0103

Compound	R.T.	QIon	Response	Conc	Unit	Qvalue
30) N-Nitrosodiphenylamine	0.00	168	0	N.D.		
31) 4-Bromophenyl-phenylether	0.00	248	0	N.D.		
32) Hexachlorobenzene	0.00	284	0	N.D.		
33) Phenanthrene	24.81	178	533	N.D.		
34) Anthracene	24.81	178	533	N.D.		
35) Di-n-butylphthalate	26.83	149	2646	N.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	32.84	228	2618	N.D.		
42) Bis(2-ethylhexyl)phthalate	33.14	149	3671	N.D.		
43) Chrysene	32.84	228	2618	N.D.		
45) Di-n-Octylphthalate	0.00	149	0	N.D.		
46) Benzo(b)fluoranthene	0.00	252	0	N.D.		
47) Benzo(k)fluoranthene	0.00	252	0	N.D.		
48) Benzo(a)pyrene	36.90	252	2260	N.D.		
49) Indeno(1,2,3-cd)pyrene	0.00	276	0	N.D.		
50) Dibenzo(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

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Quantitation Report

Data File : C:\HPCHEM\1\DATA\JAN0902\31346.D

Acq On : 11 Jan 2002 2:18 am

Sample : gcms scan 12/24

Misc :

MS Integration Params: GOOFY.P

Quant Time: Feb 5 16:40 2002

Vial: 12

Operator: 209 GALOS

Inst : GC/MS 209

Multiplr: 1.00

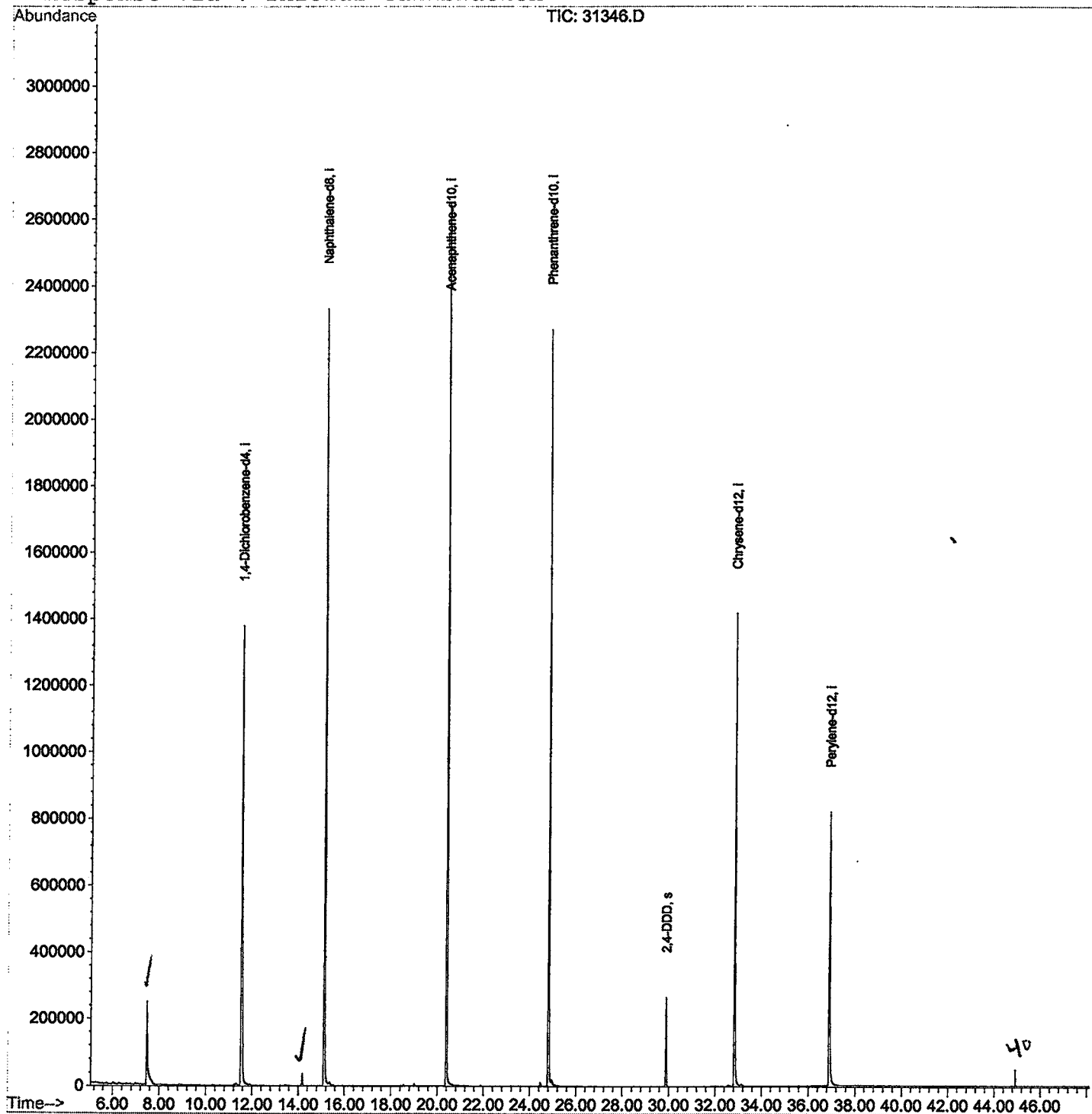
Quant Results File: GCMS0103.RES

Method : C:\HPCHEM\1\METHODS\GCMS0103.M (RTE Integrator)

Title : 209 625/TCLP calibration table

Last Update : Mon Jan 07 10:31:02 2002

Response via : Initial Calibration



LSC Area Percent Report

Data File : C:\HPCHEM\1\DATA\JAN0902\31346.D
 Acq On : 11 Jan 2002 2:18 am
 Sample : gcms scan 12/24
 Misc :
 MS Integration Params: LSCINT.P

Vial: 12
 Operator: 209 GALOS
 Inst : GC/MS 209
 Multiplr: 1.00

Method : C:\HPCHEM\1\METHODS\GCMS0103.M (RTE Integrator)
 Title : 209 625/TCLP calibration table
 Smoothing : OFF
 Sampling : 1
 Start Thrs: 0.2
 Stop Thrs : 0
 Filtering: 5
 Min Area: .1 % of largest Peak
 Max Peaks: 100
 Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >
 Peak separation: 5

Signal : TIC

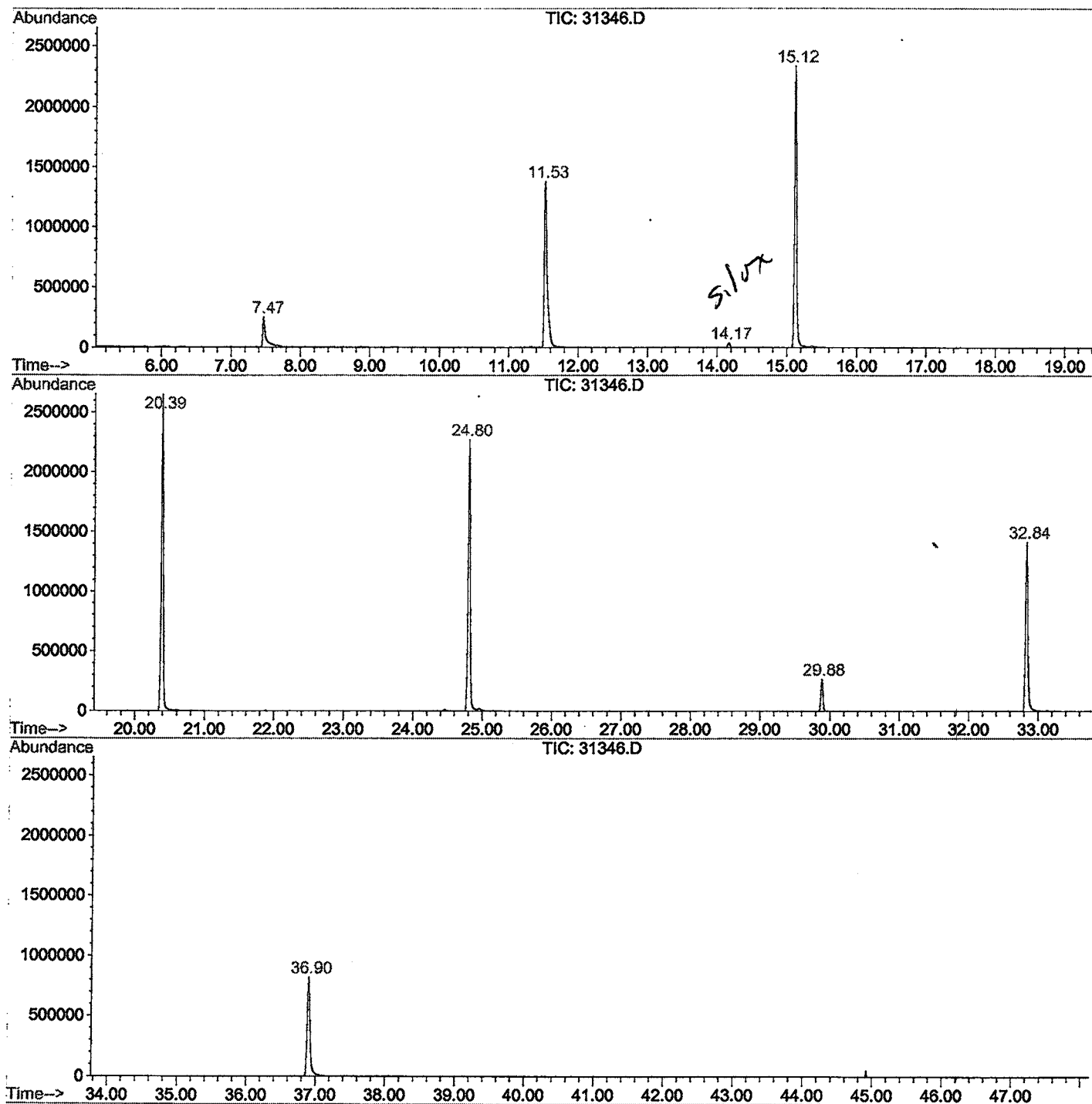
peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	peak area	peak % max.	% of total
1	7.471	260	264	283	rBV	247223	765235	14.55%	2.926%
2	11.529	701	706	727	rBV	1377089	3916765	74.47%	14.974%
3	14.173	989	994	1000	rVB2	37152	75480	1.44%	0.289%
4	15.118	1092	1097	1116	rBV	2331754	5017207	95.40%	19.181%
5	20.388	1665	1671	1690	rBV	2650934	5259228	100.00%	20.107%
6	24.803	2146	2152	2164	rBV2	2270291	4927136	93.69%	18.837%
7	29.880	2700	2705	2715	rBV	266442	542666	10.32%	2.075%
8	32.836	3021	3027	3050	rBV2	1415937	3383076	64.33%	12.934%
9	36.903	3463	3470	3488	rBV	818068	2269998	43.16%	8.678%

Sum of corrected areas: 26156791

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LSC Report - Integrated Chromatogram

File : C:\HPCHEM\1\DATA\JAN0902\31346.D
Operator : 209 GALOS
Acquired : 11 Jan 2002 2:18 am using AcqMethod GCMS0103
Instrument : GC/MS 209
Sample Name: gcms scan 12/24
Misc Info :
Vial Number: 12
Quant File :GCMS0103.RES (RTE Integrator)



31346.D GCMS0103.M

Tue Feb 05 16:58:21 2002

Library Search Compound Report

Data File : C:\HPCHEM\1\DATA\JAN0902\31346.D
Acq On : 11 Jan 2002 2:18 am
Sample : gcms scan 12/24
Misc :
MS Integration Params: LSCINT.P

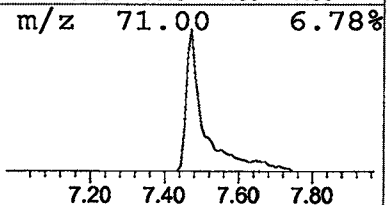
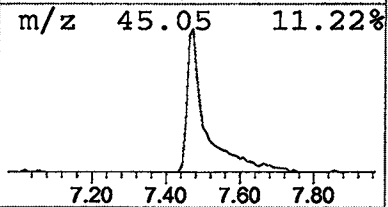
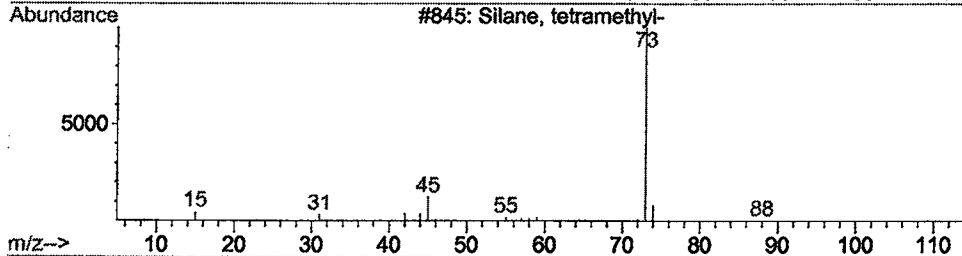
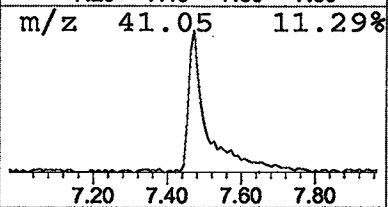
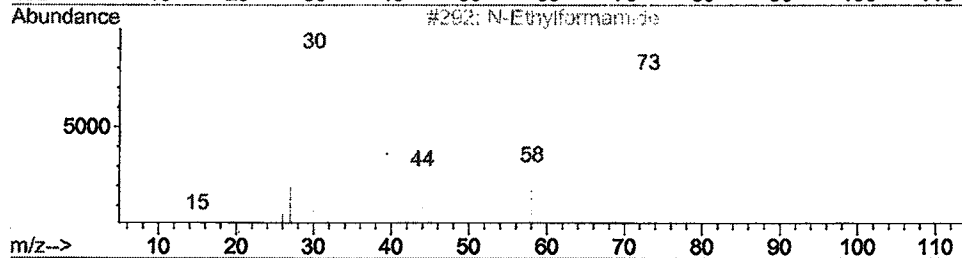
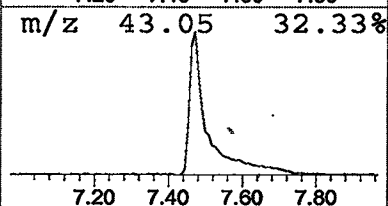
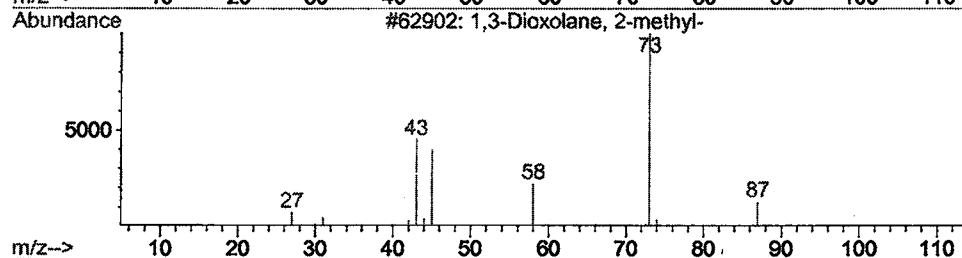
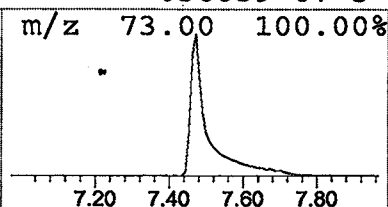
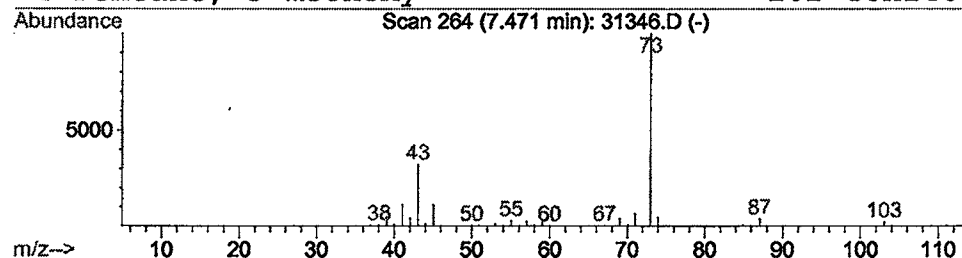
Vial: 12
Operator: 209 GALOS
Inst : GC/MS 209
Multiplr: 1.00

Quant Method : C:\HPCHEM\1\METHODS\GCMS0103.M (RTE Integrator)
Title : 209 625/TCLP calibration table
Library : C:\DATABASE\NBS75K.L

Peak Number 1 1,3-Dioxolane, 2-methyl- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.47	3.91 ug/l	765235	1,4-Dichlorobenzene-d4	11.53

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1,3-Dioxolane, 2-methyl-	88	C4H8O2	000497-26-7	9
2			N-Ethylformamide	73	C3H7NO	000627-45-2	9
3			Silane, tetramethyl-	88	C4H12Si	000075-76-3	9
4			Pentane, 3-methoxy-	102	C6H14O	036839-67-5	9



Library Search Compound Report

Data File : C:\HPCHEM\1\DATA\JAN0902\31346.D
 Acq On : 11 Jan 2002 2:18 am
 Sample : gcms scan 12/24
 Misc :
 MS Integration Params: LSCINT.P

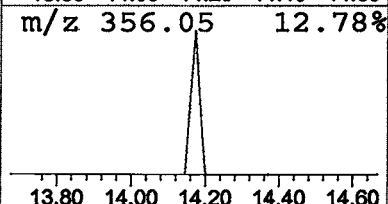
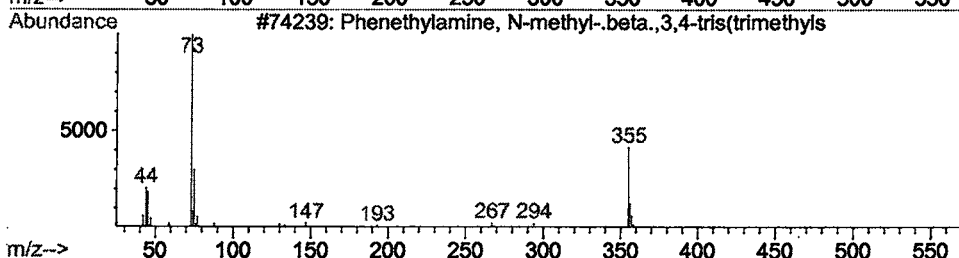
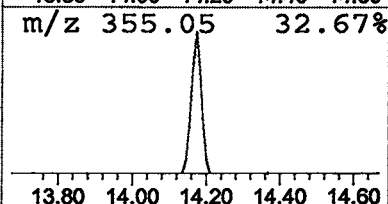
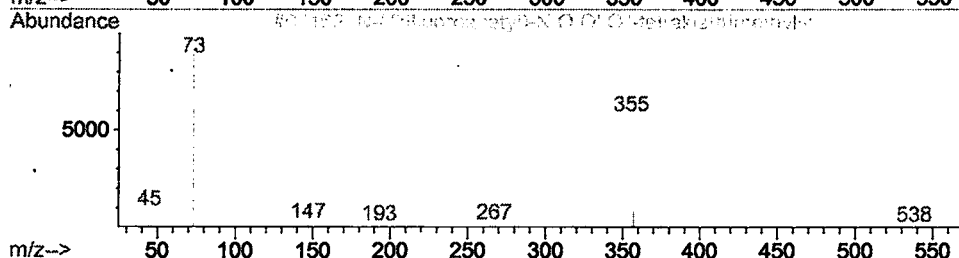
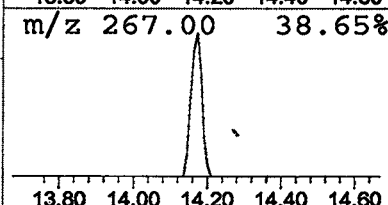
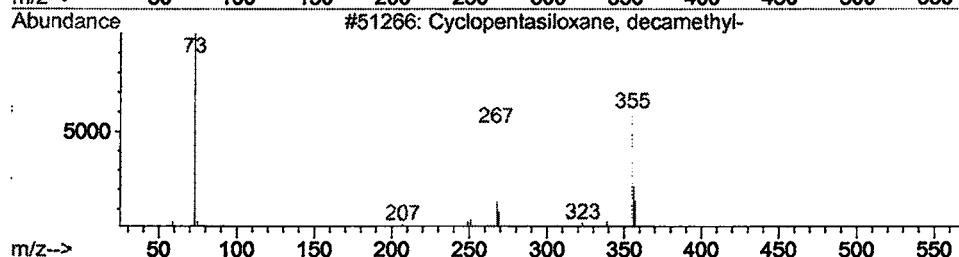
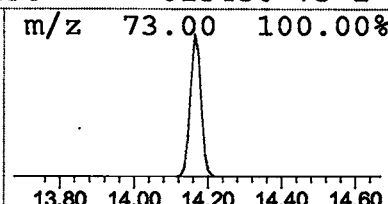
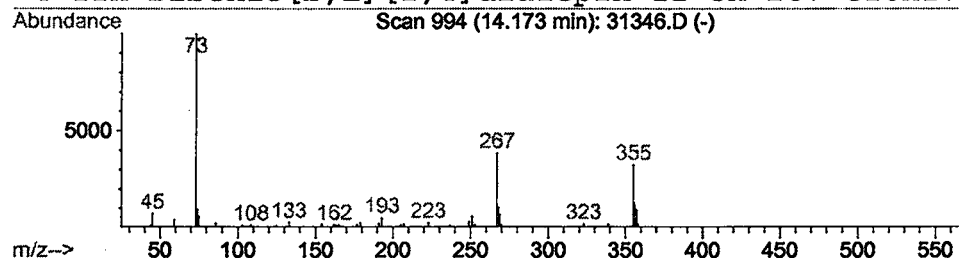
Vial: 12
 Operator: 209 GALOS
 Inst : GC/MS 209
 Multiplr: 1.00

Quant Method : C:\HPCHEM\1\METHODS\GCMS0103.M (RTE Integrator)
 Title : 209 625/TCLP calibration table
 Library : C:\DATABASE\NBS75K.L

 Peak Number 2 Cyclopentasiloxane, decamethyl Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.17	0.30 ug/l	75480	Naphthalene-d8	15.12

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Cyclopentasiloxane, decamethyl-	370	C10H30O5Si5	000541-02-6	53
2		N-(Trifluoroacetyl)-N,O,O',O''-tetr	553	C22H42F3NO4Si4	000000-00-0	35
3		Phenethylamine, N-methyl-.beta.,3,4	399	C18H37NO3Si3	010538-85-9	32
4		11H-Dibenzo[b,E][1,4]diazepin-11-on	267	C16H17N3O	013450-73-2	27



Tentatively Identified Compound (LSC) summary

Operator ID: 209 GALOS Date Acquired: 11 Jan 2002 2:18 am
 Data File: C:\HPCHEM\1\DATA\JAN0902\31346.D
 Name: gcms scan 12/24
 Misc:
 Method: C:\HPCHEM\1\METHODS\GCMS0103.M (RTE Integrator)
 Title: 209 625/TCLP calibration table
 Library Searched: C:\DATABASE\NBS75K.L

TIC Top Hit name	RT	EstConc	Units	Area	IntStd	ISRT	ISArea	ISConc
1,3-Dioxolane, 2-met	7.47	3.9	ug/l	765235	ISTD01	11.53	3916770	20.0
Cyclopentasiloxane,	14.17	0.3	ug/l	75480	ISTD02	15.12	5017210	20.0

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