

User: Vinci, David L
 Westchester Dept. of Labs & Research
 Date: 03/05/2002 Time: 12:27:02

Sample: AE02721
 Analysis: \$GCMS GCMS Scan For Unknowns
 Units: ug
 Started: 3/5/02 12:26 Ended: 3/5/02 12:26 Analyst: DLV
 Page: 1

	Component Name	Viol	Result	Qualifier	MDL
1	Other unknown compounds		.		
2	Surr_2,4-DDD		77 ✓		5.0

Comments for Sample AE02721 - Analysis \$GCMS

Westchester Dept. of Labs & Research

User: Vinci, David L.

Date: 03-05-2002 Time: 12:27:11

The GCMS scan, from 35 to 550 amu, does not reveal the presence of unusual compounds. The recoveries for 14 of the 43 compounds in the LCS Standard were above their acceptance ranges.

Quantitation Report

(Not Reviewed)

Data File : C:\MSDCHEM\1\5973N\02FEB28\02721.D

Vial: 6

Acq On : 28 Feb 2002 6:50 pm

Operator:

Sample : 508 scan

Inst : Instrumen

Misc : February 28, 2002

Multiplr: 1.00

MS Integration Params: SPLIT.P

Quant Time: Mar 01 09:03:02 2002

Quant Results File: HP022102.RES

Quant Method : C:\MSDCHEM\1...\HP022102.M (RTE Integrator)

Title : 508 scan method

Last Update : Mon Feb 25 10:47:59 2002

Response via : Initial Calibration

DataAcq Meth : HP020702

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	9.72	152	1427827	20.00	ug/L	0.00
10) Naphthalene-d8	13.20	136	6173848	20.00	ug/L	0.00
17) Acenaphthene-d10	18.33	164	3085827	20.00	ug/L	0.00
29) Phenanthrene-d10	22.63	188	5494854	20.00	ug/L	-0.01
38) Chrysene-d12	30.49	240	4063252	20.00	ug/L	-0.03
44) Perylene-d12	34.41	264	2663213	20.00	ug/L	-0.04

System Monitoring Compounds

37) 2,4-DDD	27.72	235	282354	3.85	ug/L	0.00
Spiked Amount	5.000		Recovery	=	77.00%	

Target Compounds

					Qvalue
20) Dimethylphthalate	18.34	163	501807	2.68 ug/L	# 57
21) 2,6-Dinitrotoluene	18.33	165	396339	9.17 ug/L	# 27

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

02721.D HP022102.M Fri Mar 01 09:03:03 2002

Page 1

Quantitation Report

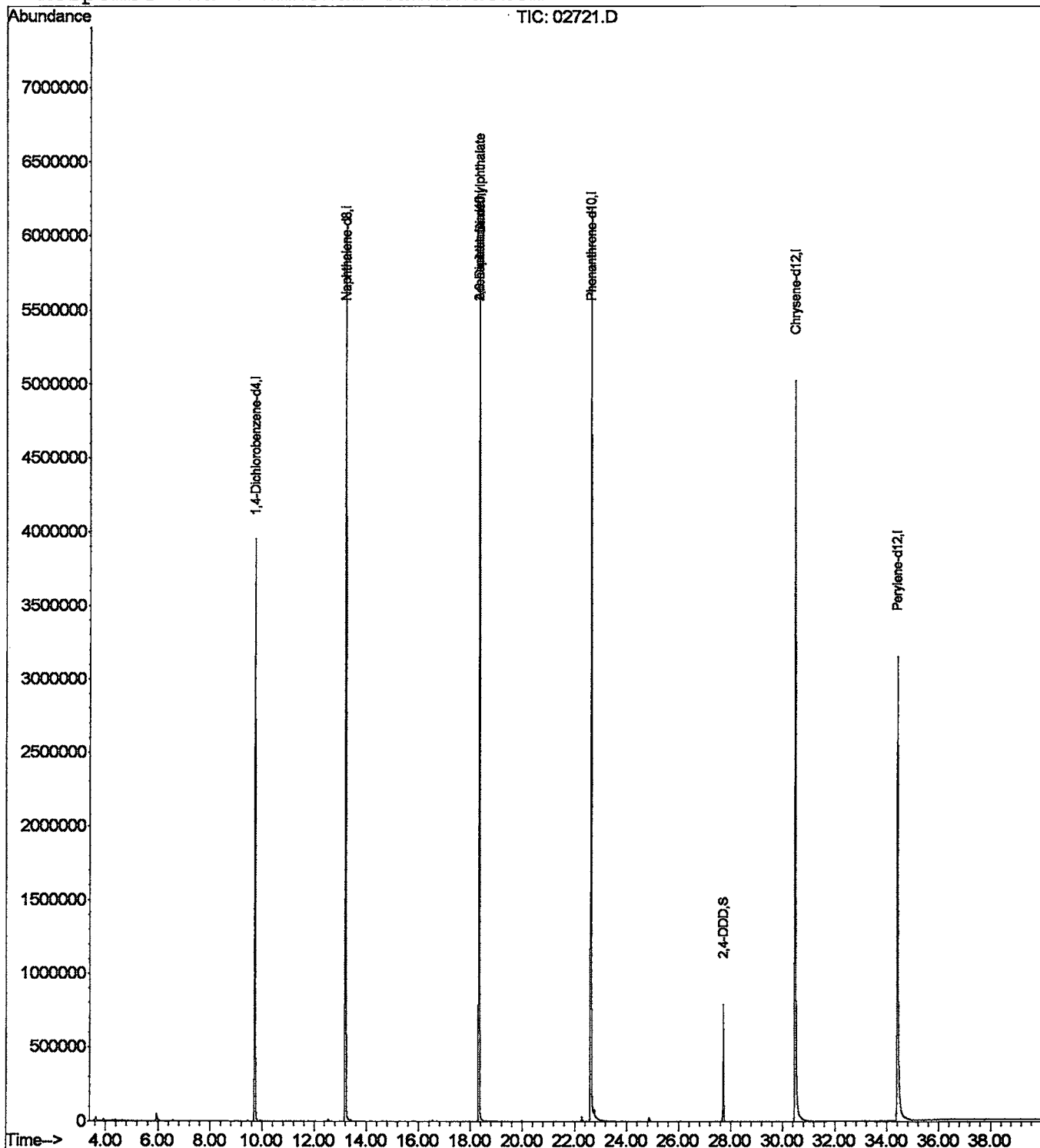
(Not Reviewed)

Data File : C:\MSDCHEM\1\5973N\02FEB28\02721.D
 Acq On : 28 Feb 2002 6:50 pm
 Sample : 508 scan
 Misc : February 28, 2002
 MS Integration Params: SPLIT.P
 Quant Time: Mar 1 9:03 2002

Vial: 6
 Operator:
 Inst : Instrumen
 Multiplr: 1.00

Quant Results File: HP022102.R

Method : C:\MSDCHEM\1\METHODS\METHODS\HP022102.M (RTE Integrator)
 Title : 508 scan method
 Last Update : Mon Feb 25 10:47:59 2002
 Response via : Initial Calibration



LSC Area Percent Report

Data File : C:\MSDCHEM\1\5973N\02FEB28\02721.D
 Acq On : 28 Feb 2002 6:50 pm
 Sample : 508 scan
 Misc : February 28, 2002
 MS Integration Params: LSCINT.P

Vial: 6
 Operator:
 Inst : Instrumen
 Multiplr: 1.00

Method : C:\MSDCHEM\1\METHODS\METHODS\HP022102.M (RTE Integrator)
 Title : 508 scan method
 Smoothing : OFF
 Sampling : 2
 Start Thrs: 0.2
 Stop Thrs : 0

Filtering: 5
 Min Area: 100 Area counts
 Max Peaks: 20
 Peak Location: TOP

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

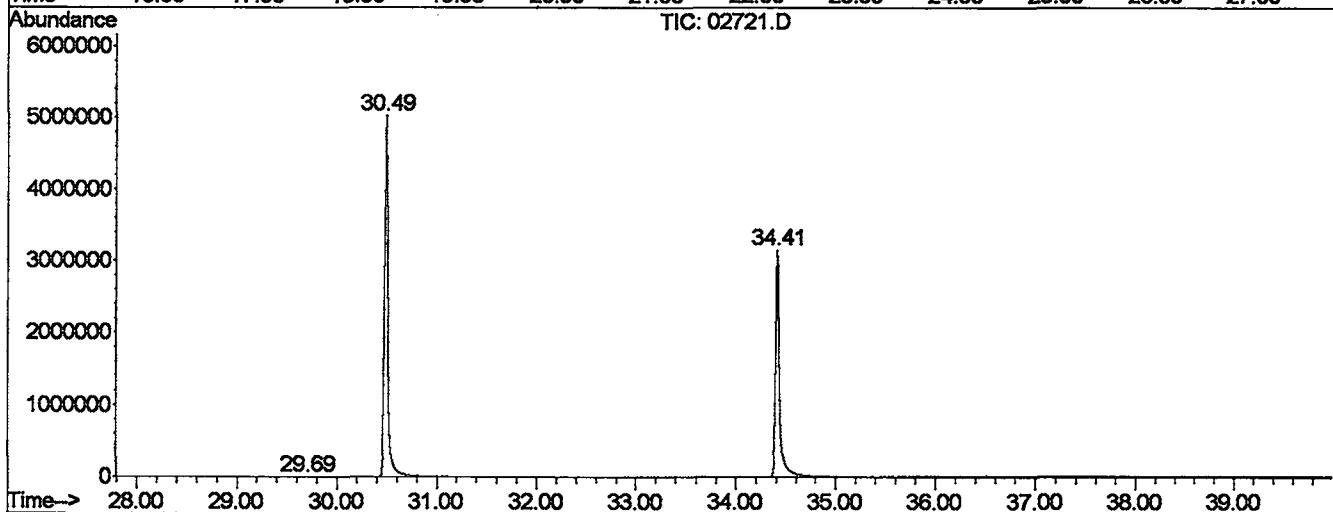
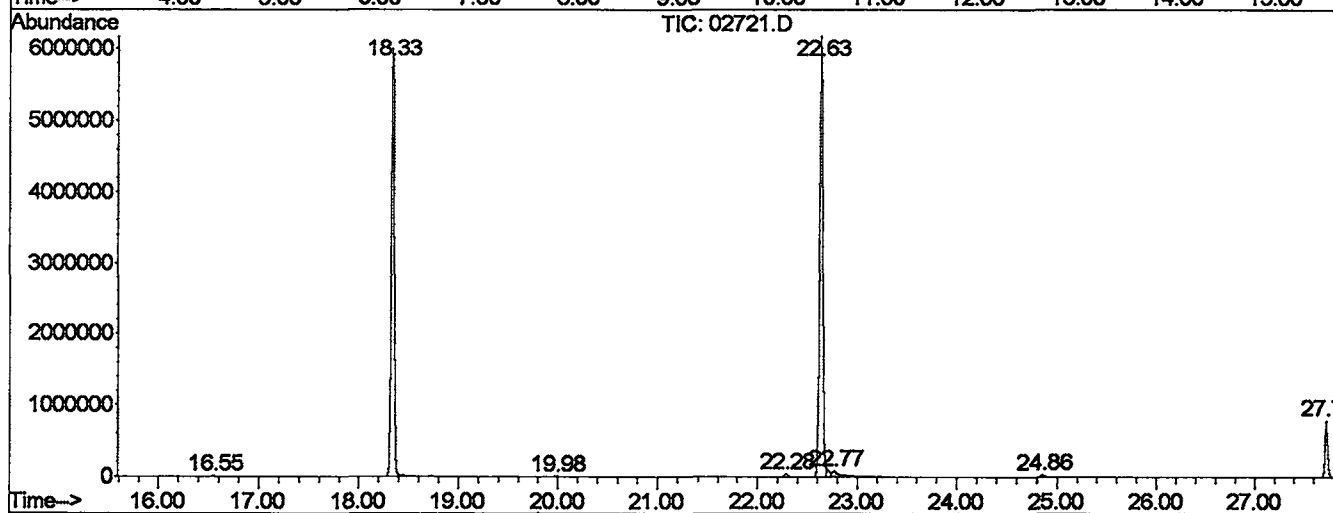
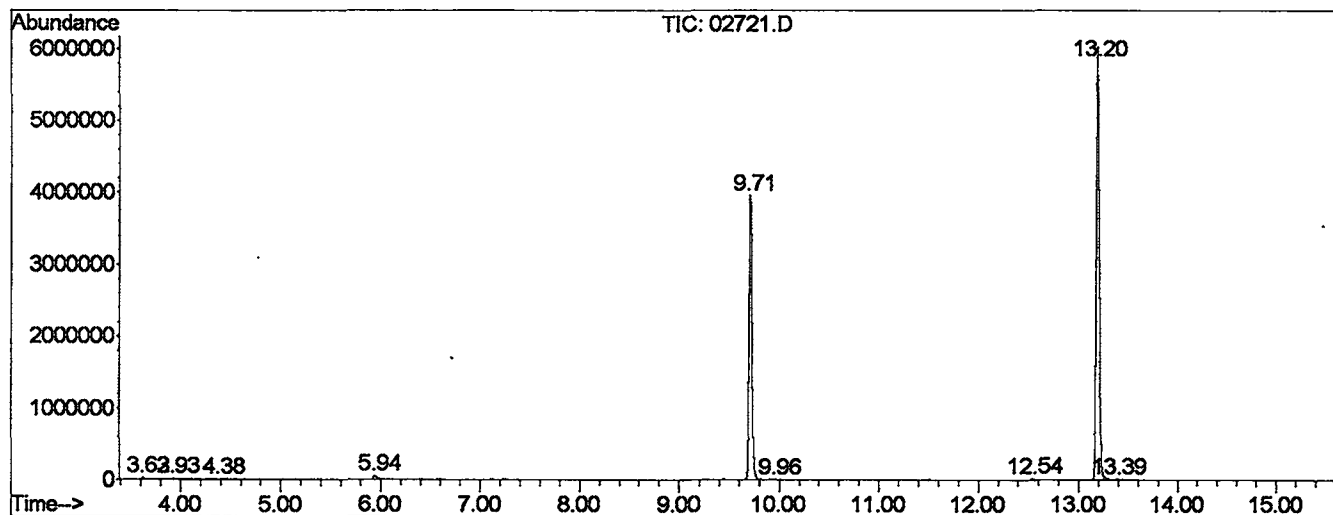
Signal : TIC

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.622	19	22	29	rBV2	27309	52411	0.39%	0.076%
2	3.927	45	49	55	rVB	14335	22817	0.17%	0.033%
3	4.379	85	89	95	rBV2	8389	18429	0.14%	0.027%
4	5.936	225	227	241	rBV	51620	134636	1.00%	0.195%
5	9.707	557	561	567	rBV	3950701	7966835	58.90%	11.521%
6	9.955	579	583	591	rVB5	5731	19370	0.14%	0.028%
7	12.540	807	812	817	rBV	16888	26953	0.20%	0.039%
8	13.195	865	870	875	rBV	6028170	11618641	85.90%	16.802%
9	13.387	883	887	897	rVB2	11957	38017	0.28%	0.055%
10	16.548	1159	1167	1171	rBV2	8802	19813	0.15%	0.029%
11	18.332	1319	1325	1331	rBV	6002124	12647942	93.51%	18.290%
12	19.980	1467	1471	1475	rBV2	5956	17003	0.13%	0.025%
13	22.283	1671	1675	1681	rBV2	35670	76233	0.56%	0.110%
14	22.633	1701	1706	1711	rBV2	6175897	13526062	100.00%	19.560%
15	22.768	1715	1718	1733	rVB2	72713	301511	2.23%	0.436%
16	24.857	1899	1903	1915	rBV	26462	66389	0.49%	0.096%
17	27.724	2153	2157	2163	rBV	791102	1363685	10.08%	1.972%
18	29.688	2325	2331	2345	rBV2	2398	15726	0.12%	0.023%
19	30.490	2397	2402	2449	rVV	5022547	12059650	89.16%	17.440%
20	34.407	2743	2749	2799	rBV	3144609	9158962	67.71%	13.245%

Sum of corrected areas: 69151085

LSC Report - Integrated Chromatogram

File : C:\MSDCHEM\1\5973N\02FEB28\02721.D
 Operator :
 Acquired : 28 Feb 2002 6:50 pm using AcqMethod HP020702
 Instrument : Instrumen
 Sample Name: 508 scan
 Misc Info : February 28, 2002
 Vial Number: 6
 Quant File :HP022102.RES (RTE Integrator)



Library Search Compound Report

Data File : C:\MSDCHEM\1\5973N\02FEB28\02721.D
Acq On : 28 Feb 2002 6:50 pm
Sample : 508 scan
Misc : February 28, 2002
MS Integration Params: LSCINT.P

Vial: 6
Operator:
Inst : Instrumen
Multiplr: 1.00

Quant Method : C:\MSDCHEM\1...\HP022102.M (RTE Integrator)

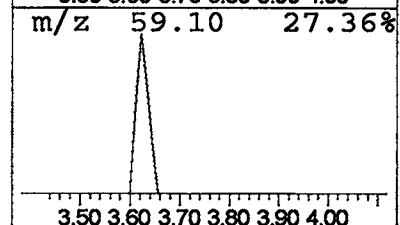
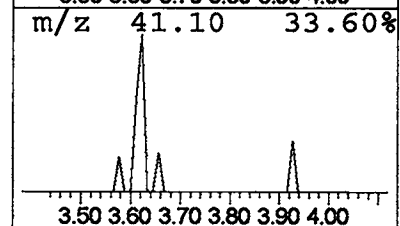
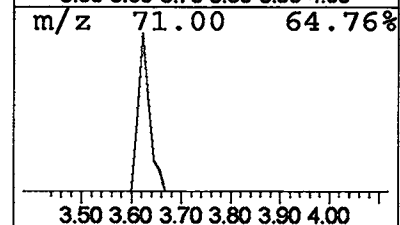
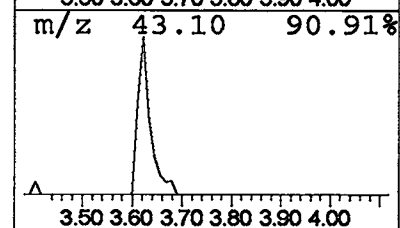
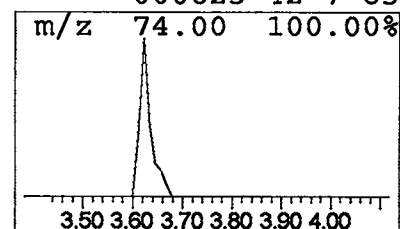
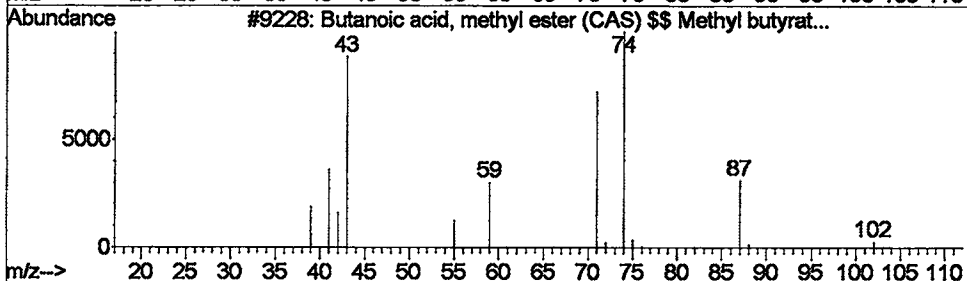
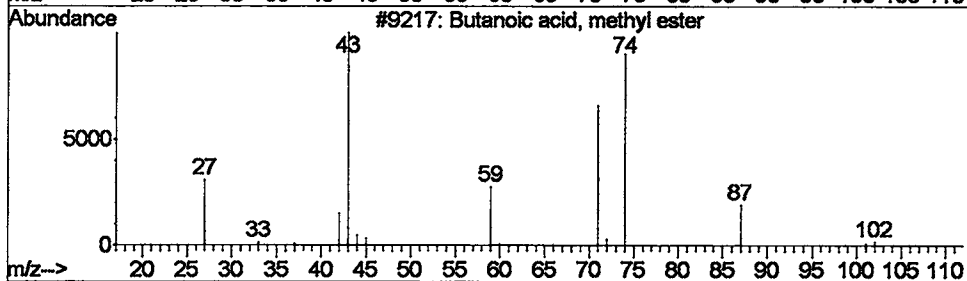
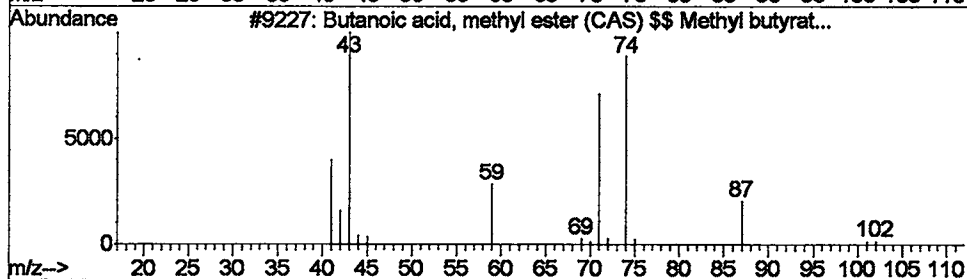
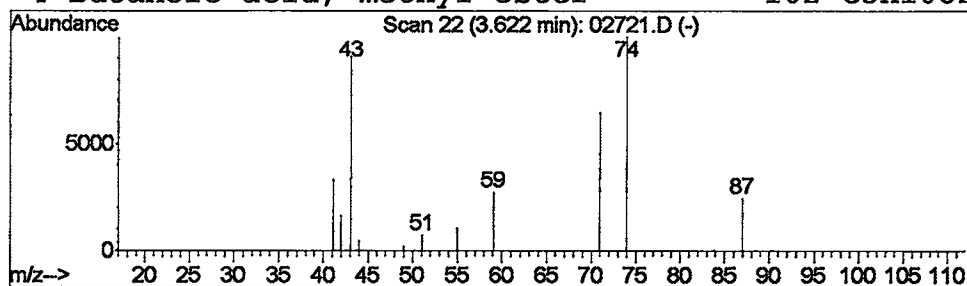
Title : 508 scan method

Library : C:\DATABASE\WILEY7N.L

Peak Number 1 Butanoic acid, methyl ester... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.62	0.13 ug/L	52411	1,4-Dichlorobenzene-d4	9.72

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Butanoic acid, methyl ester (CAS...	102	C5H10O2	000623-42-7	83
2			Butanoic acid, methyl ester	102	C5H10O2	000623-42-7	83
3			Butanoic acid, methyl ester (CAS...	102	C5H10O2	000623-42-7	83
4			Butanoic acid, methyl ester	102	C5H10O2	000623-42-7	83



Library Search Compound Report

Data File : C:\MSDCHEM\1\5973N\02FEB28\02721.D
 Acq On : 28 Feb 2002 6:50 pm
 Sample : 508 scan
 Misc : February 28, 2002
 MS Integration Params: LSCINT.P

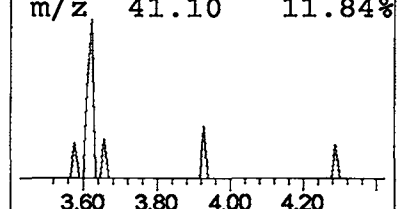
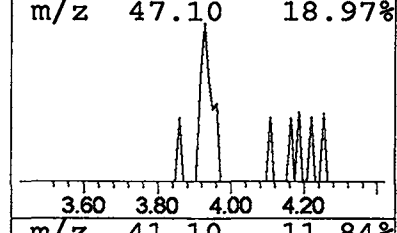
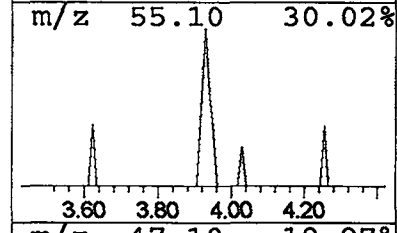
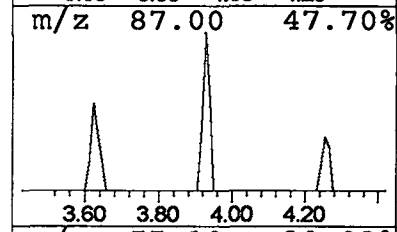
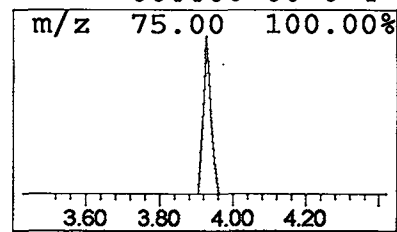
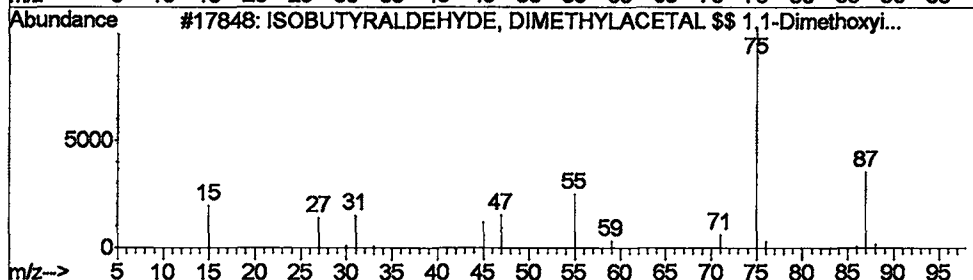
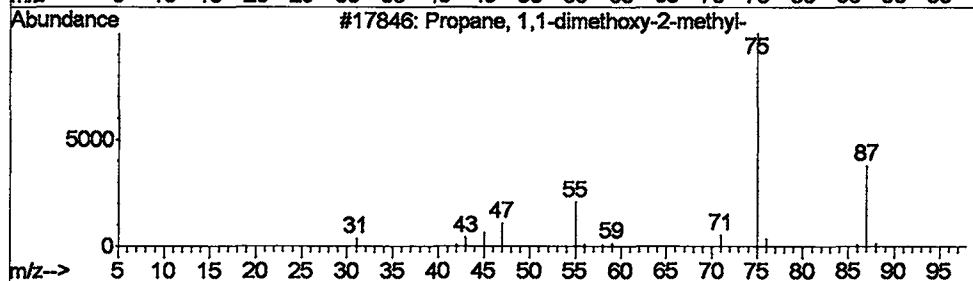
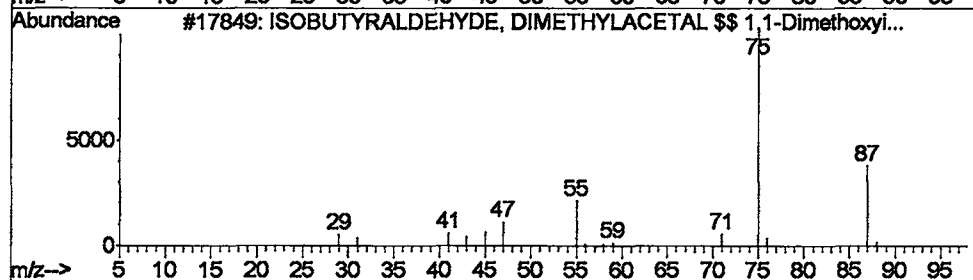
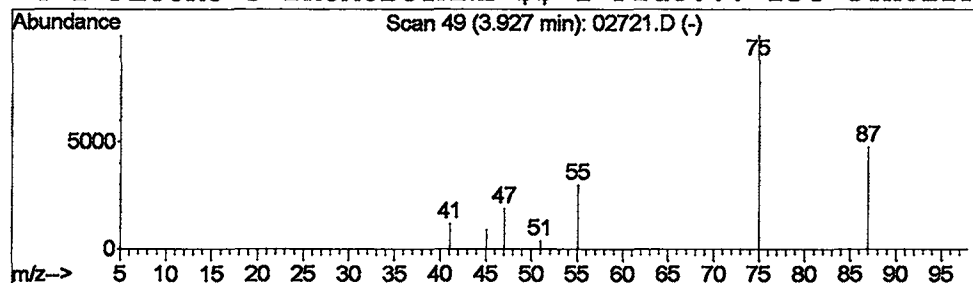
Vial: 6
 Operator:
 Inst : Instrumen
 Multiplr: 1.00

Quant Method : C:\MSDCHEM\1...\HP022102.M (RTE Integrator)
 Title : 508 scan method
 Library : C:\DATABASE\WILEY7N.L

 Peak Number 2 ISOBUTYRALDEHYDE, DIMETHYLA... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.93	0.06 ug/L	22817	1,4-Dichlorobenzene-d4	9.72

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		ISOBUTYRALDEHYDE, DIMETHYLACETAL...	118	C6H14O2	000000-00-0	9
2		Propane, 1,1-dimethoxy-2-methyl-	118	C6H14O2	041632-89-7	9
3		ISOBUTYRALDEHYDE, DIMETHYLACETAL...	118	C6H14O2	000000-00-0	9
4		1-FLUORO-3-BROMOBUTANE \$\$ 1-Fluo...	154	C4H8BrF	000000-00-0	4



Library Search Compound Report

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Acq On : 28 Feb 2002 6:50 pm
Sample : 508 scan
Misc : February 28, 2002
MS Integration Params: LSCINT.P

Vial: 6
Operator:
Inst : Instrumen
Multiplr: 1.00

Quant Method : C:\MSDCHEM\1...\HP022102.M (RTE Integrator)

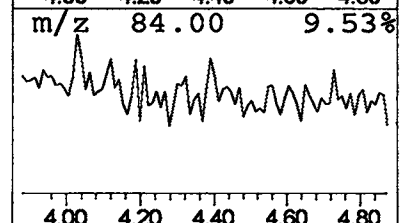
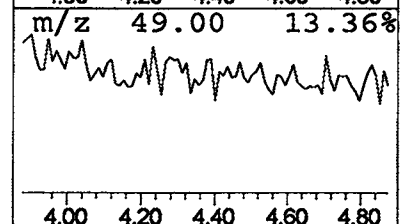
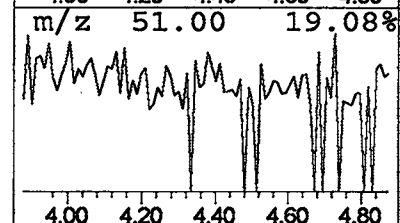
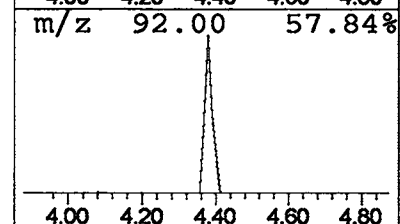
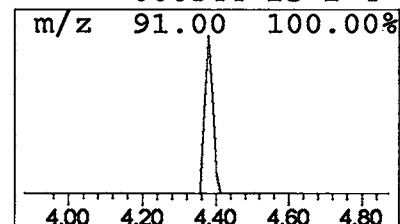
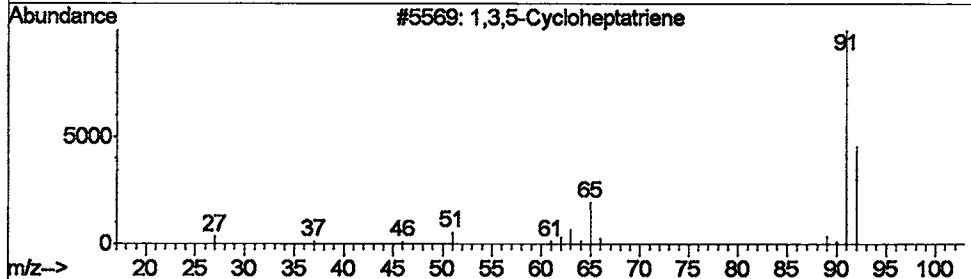
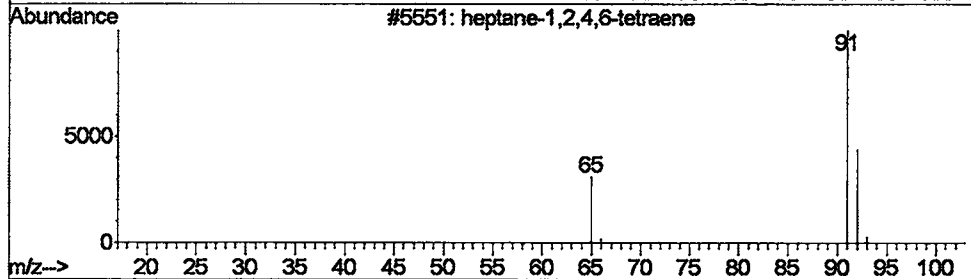
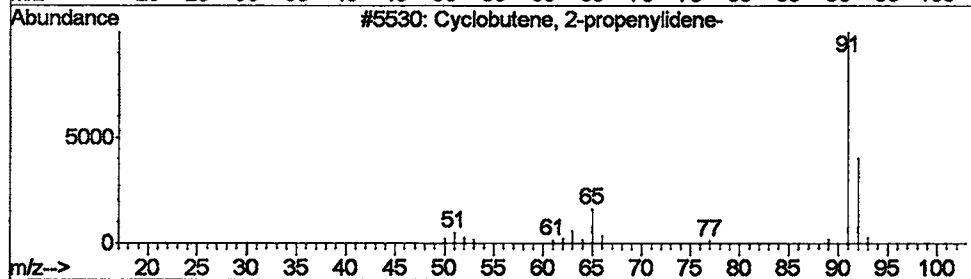
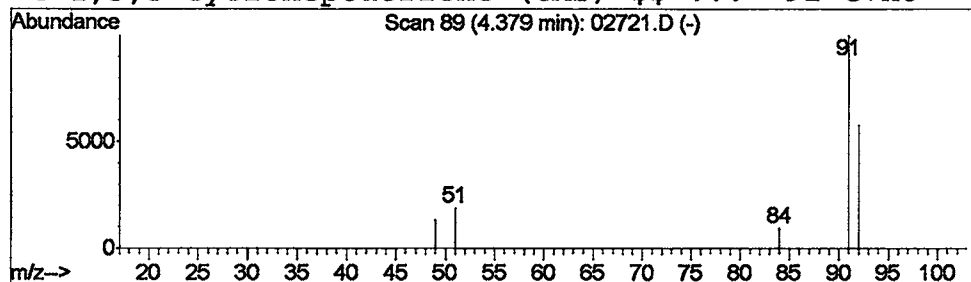
Title : 508 scan method

Library : C:\DATABASE\WILEY7N.L

Peak Number 3 Cyclobutene, 2-propenylidene- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.38	0.05 ug/L	18429	1,4-Dichlorobenzene-d4	9.72

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclobutene, 2-propenylidene-	92	C7H8	052097-85-5	4
2			heptane-1,2,4,6-tetraene	92	C7H8	069961-24-6	4
3			1,3,5-Cycloheptatriene	92	C7H8	000544-25-2	4
4			1,3,5-Cycloheptatriene (CAS) \$\$...	92	C7H8	000544-25-2	4



Library Search Compound Report

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 Sample : 508 scan
 Misc : February 28, 2002
 MS Integration Params: LSCINT.P

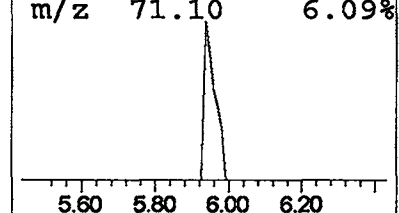
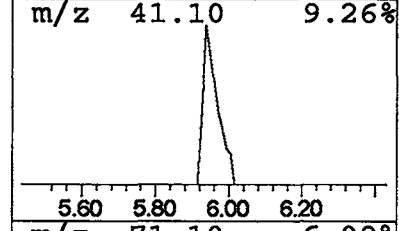
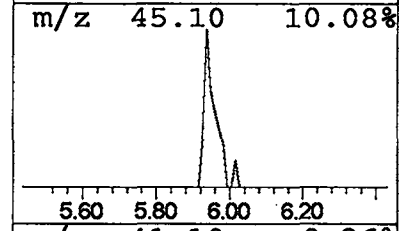
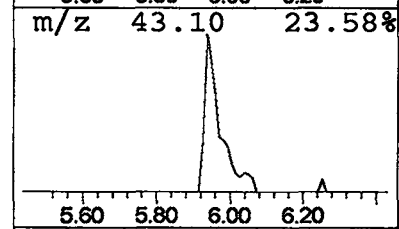
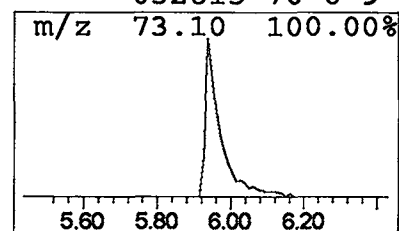
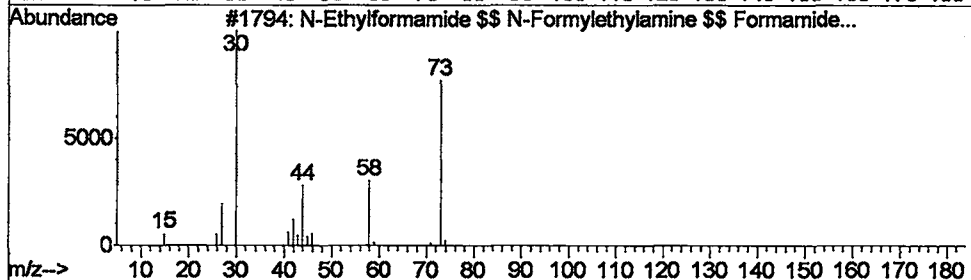
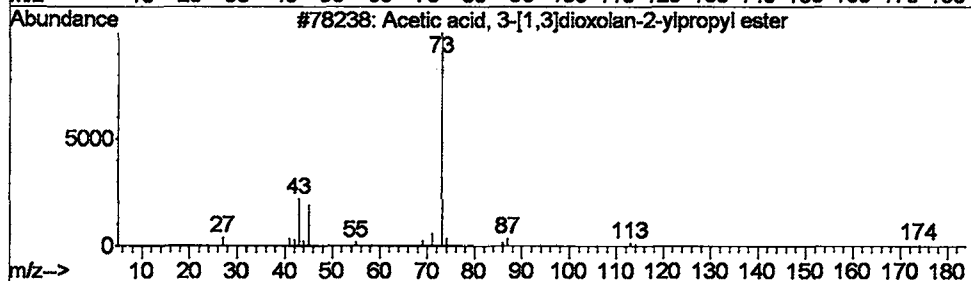
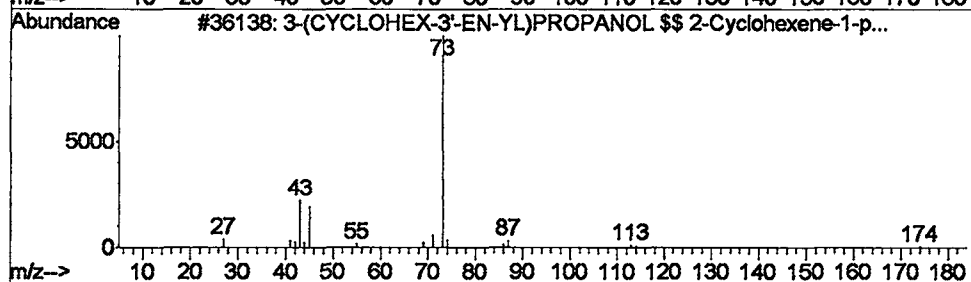
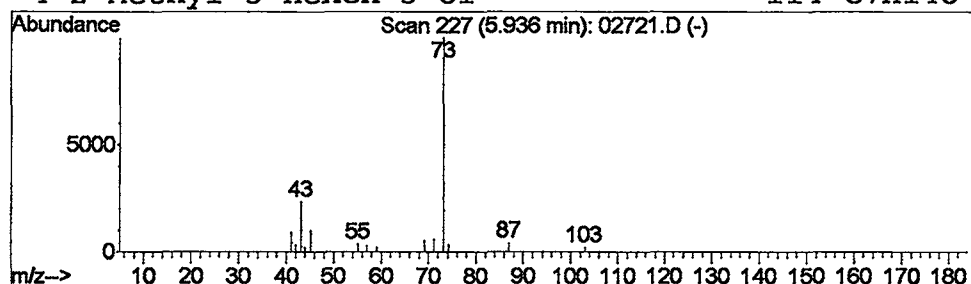
Vial: 6
 Operator:
 Inst : Instrumen
 Multiplr: 1.00

Quant Method : C:\MSDCHEM\1...\HP022102.M (RTE Integrator)
 Title : 508 scan method
 Library : C:\DATABASE\WILEY7N.L

 Peak Number 4 3-(CYCLOHEX-3'-EN-YL)PROPAN... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.94	0.34 ug/L	134636	1,4-Dichlorobenzene-d4	9.72

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		3-(CYCLOHEX-3'-EN-YL)PROPANOL \$\$...	174	C8H14O4	015745-87-6	36
2		Acetic acid, 3-[1,3]dioxolan-2-y...	174	C8H14O4	000000-00-0	36
3		N-Ethylformamide \$\$ N-Formylethy...	73	C3H7NO	000627-45-2	9
4		2-Methyl-5-hexen-3-ol	114	C7H14O	032815-70-6	9



Library Search Compound Report

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 Acq On : 28 Feb 2002 6:50 pm
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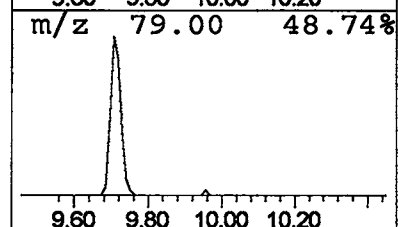
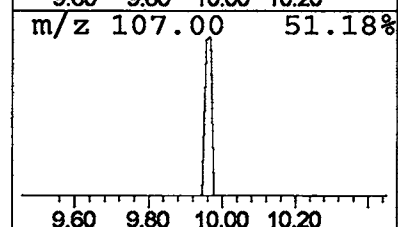
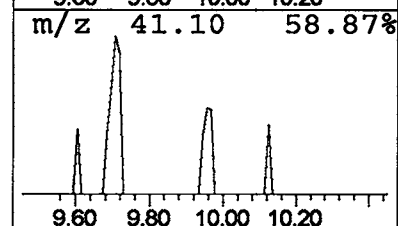
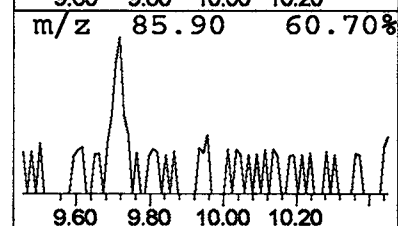
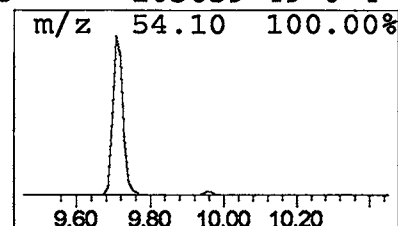
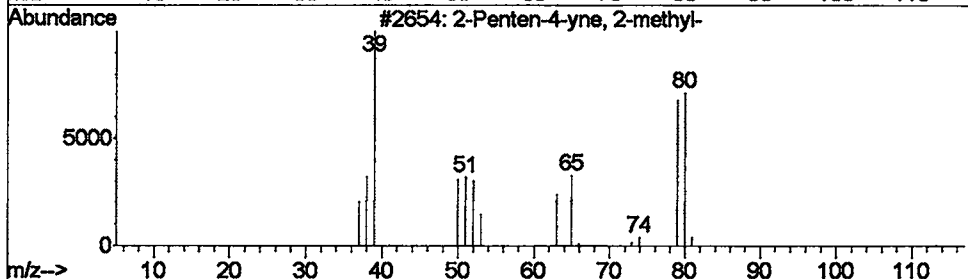
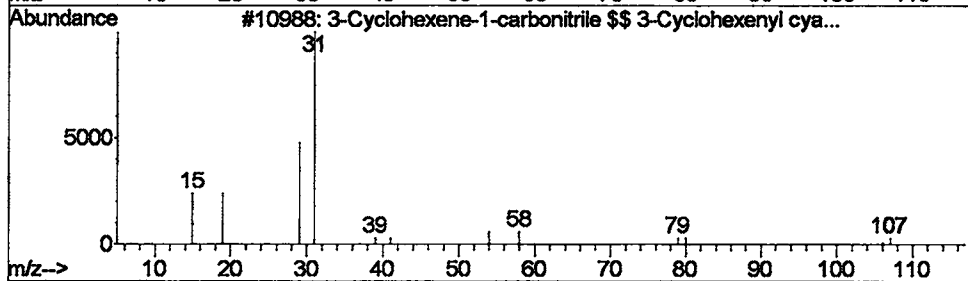
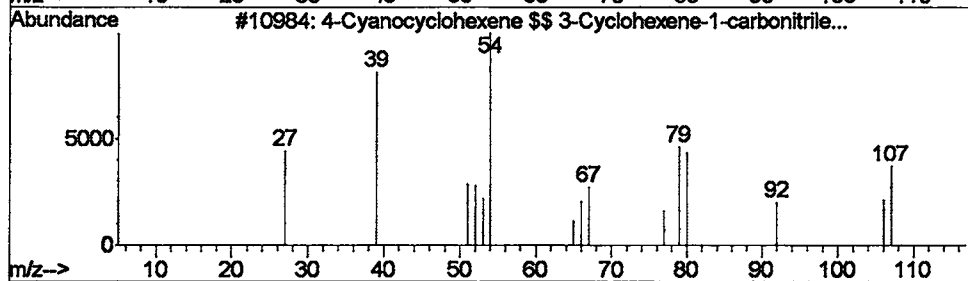
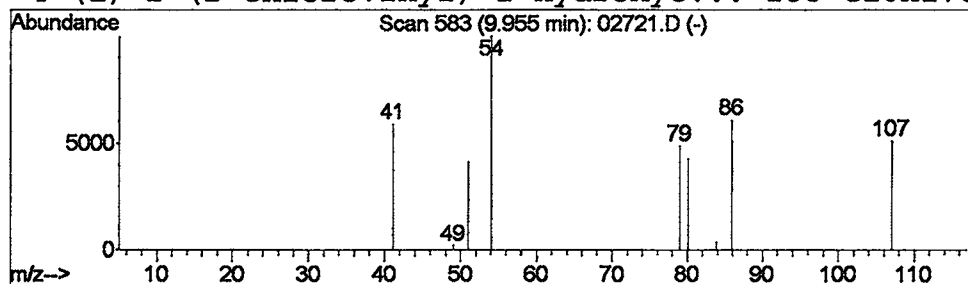
Vial: 6
 Operator:
 Inst : Instrumen
 Multiplr: 1.00

Quant Method : C:\MSDCHEM\1...\HP022102.M (RTE Integrator)
 Title : 508 scan method
 Library : C:\DATABASE\WILEY7N.L

 Peak Number 5 4-Cyanocyclohexene \$\$ 3-Cyc... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.96	0.05 ug/L	19370	1,4-Dichlorobenzene-d4	9.72

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	4-Cyanocyclohexene \$\$ 3-Cyclohex...	107	C7H9N	000100-45-8	86	
2	3-Cyclohexene-1-carbonitrile \$\$...	107	C7H9N	000100-45-8	9	
3	2-Penten-4-yne, 2-methyl-	80	C6H8	000000-00-0	4	
4	(Z)-2-(2-chlorovinyl)-1-hydroxyc...	188	C10H17ClO	103659-49-0	4	



Library Search Compound Report

Data File : C:\MSDCHEM\1\5973N\02FEB28\02721.D
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Sample : 508 scan
Misc : February 28, 2002
MS Integration Params: LSCINT.P

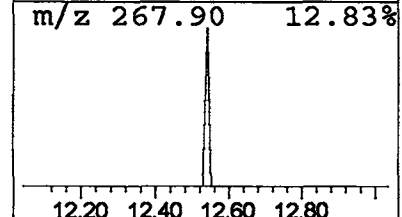
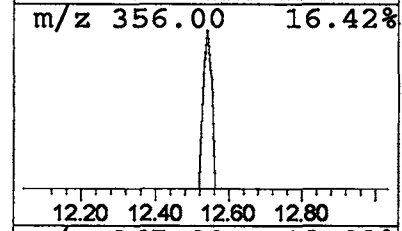
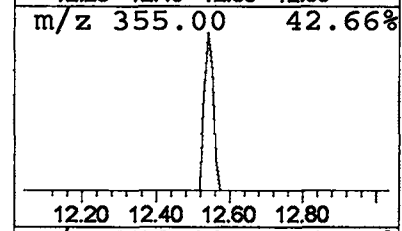
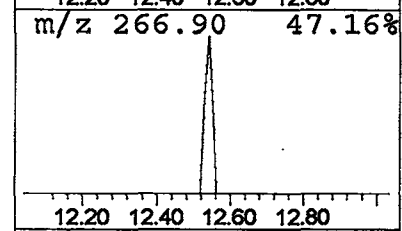
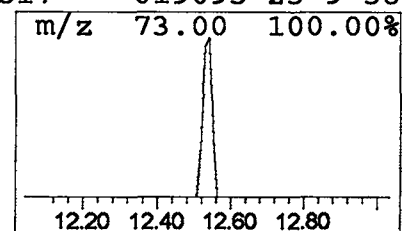
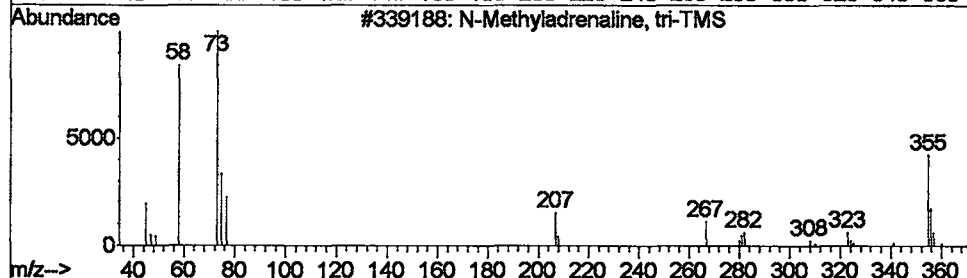
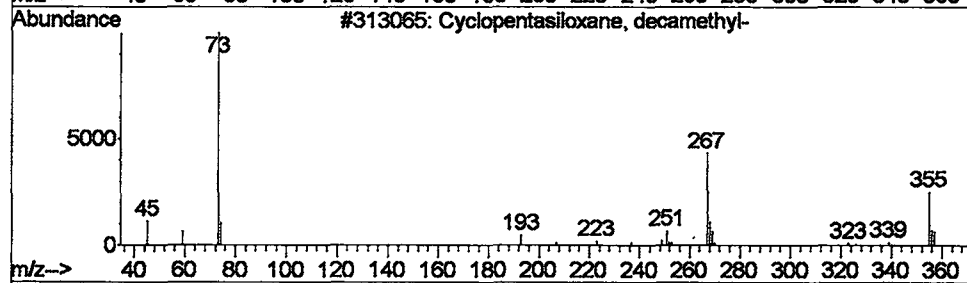
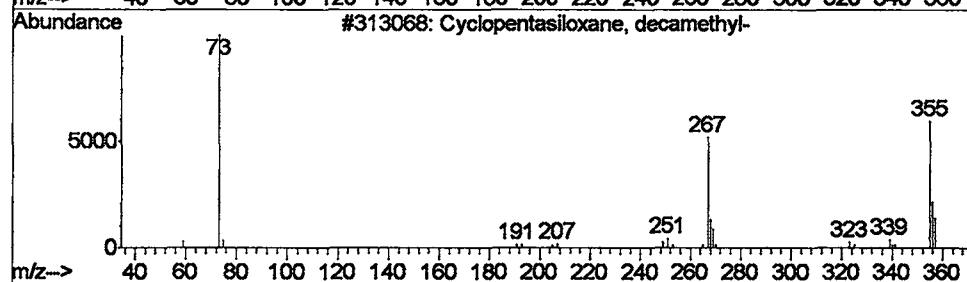
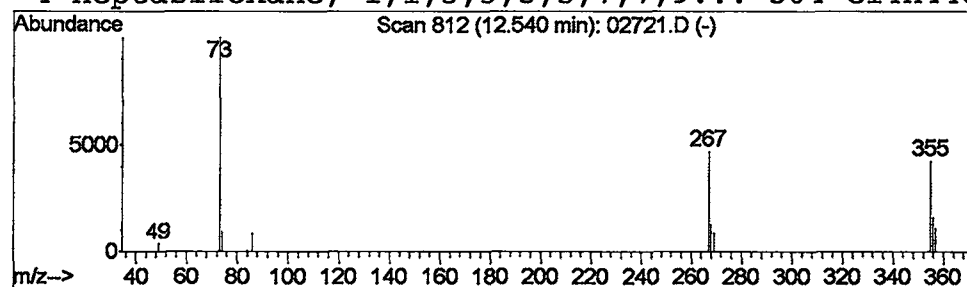
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Operator:
Inst : Instrumen
Multiplr: 1.00

Quant Method : C:\MSDCHEM\1...\HP022102.M (RTE Integrator)
Title : 508 scan method
Library : C:\DATABASE\WILEY7N.L

Peak Number 6 Cyclopentasiloxane, decamet... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.54	0.05 ug/L	26953	Naphthalene-d8	13.20

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Cyclopentasiloxane, decamethyl-	370	C10H30O5Si5	000541-02-6	83
2		Cyclopentasiloxane, decamethyl-	370	C10H30O5Si5	000541-02-6	64
3		N-Methyladrenaline, tri-TMS	413	C19H39NO3Si3	000000-00-0	40
4		Heptasiloxane, 1,1,3,3,5,5,7,7,9...	504	C14H44O6Si7	019095-23-9	38



Library Search Compound Report

Data File : C:\MSDCHEM\1\5973N\02FEB28\02721.D
Acq On : 28 Feb 2002 6:50 pm
Sample : 508 scan
Misc : February 28, 2002
MS Integration Params: LSCINT.P

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Operator:
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Multiplr: 1.00

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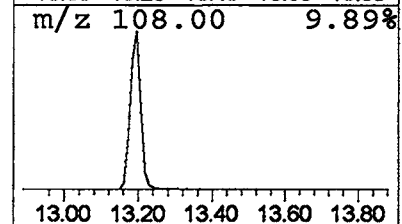
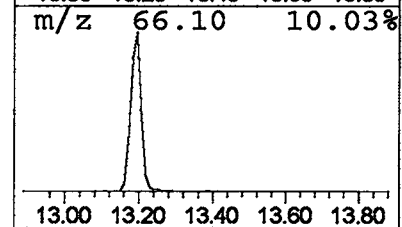
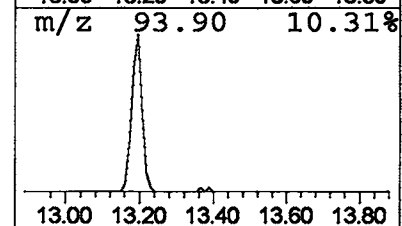
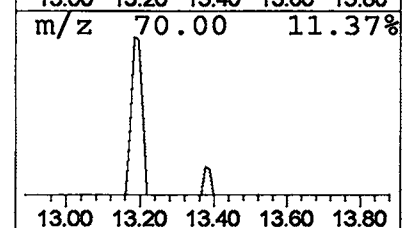
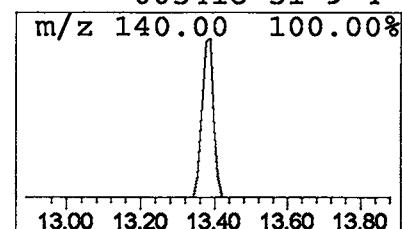
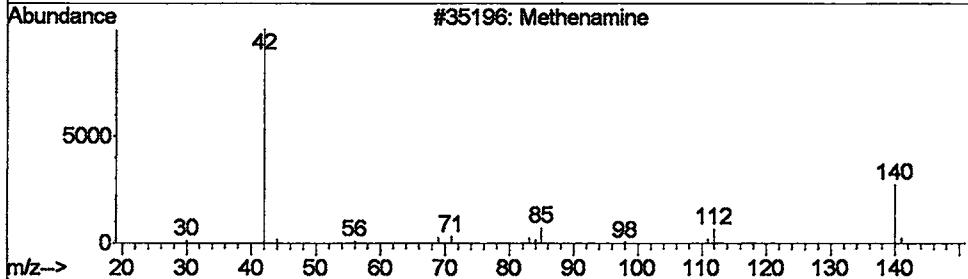
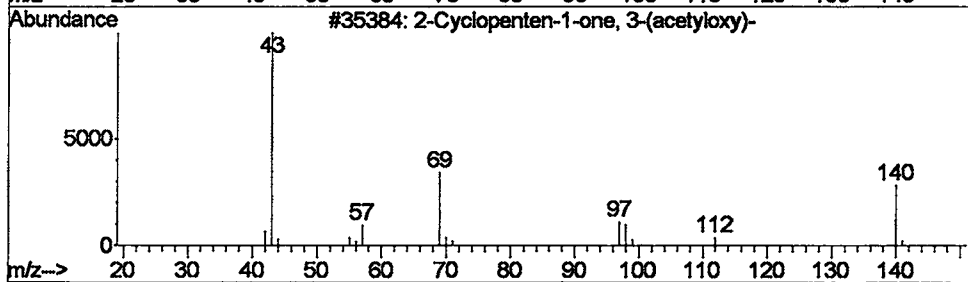
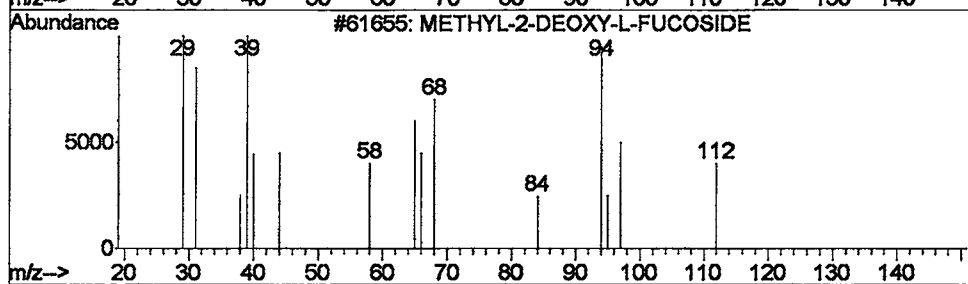
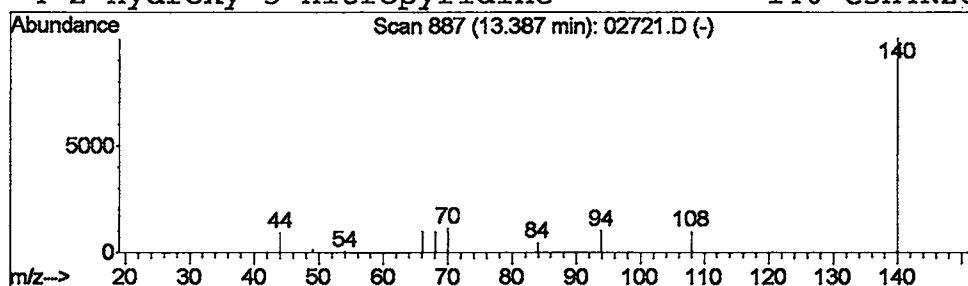
Title : 508 scan method

Library : C:\DATABASE\WILEY7N.L

Peak Number 7 METHYL-2-DEOXY-L-FUCOSIDE Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.39	0.07 ug/L	38017	Naphthalene-d8	13.20

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		METHYL-2-DEOXY-L-FUCOSIDE	162	C7H14O4	000000-00-0	9
2		2-Cyclopenten-1-one, 3-(acetyloxy)-	140	C7H8O3	018766-96-6	5
3		Methenamine	140	C6H12N4	000100-97-0	4
4		2-Hydroxy-5-nitropyridine	140	C5H4N2O3	005418-51-9	4



Library Search Compound Report

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Sample : 508 scan
Misc : February 28, 2002
MS Integration Params: LSCINT.P

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Operator:
Inst : Instrumen
Multiplr: 1.00

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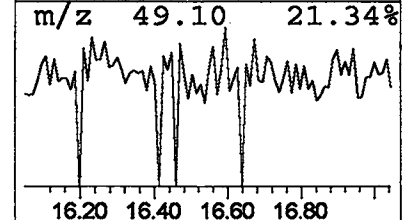
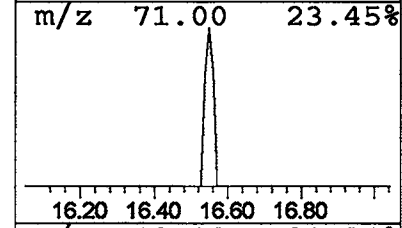
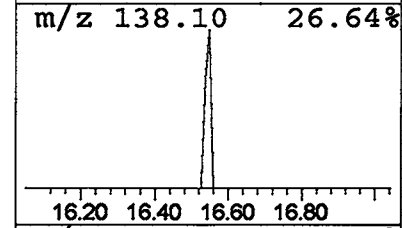
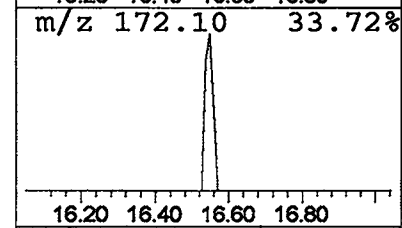
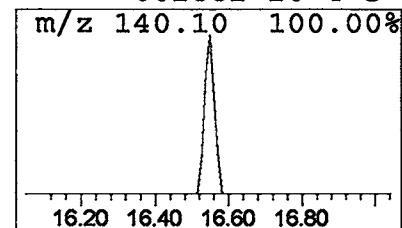
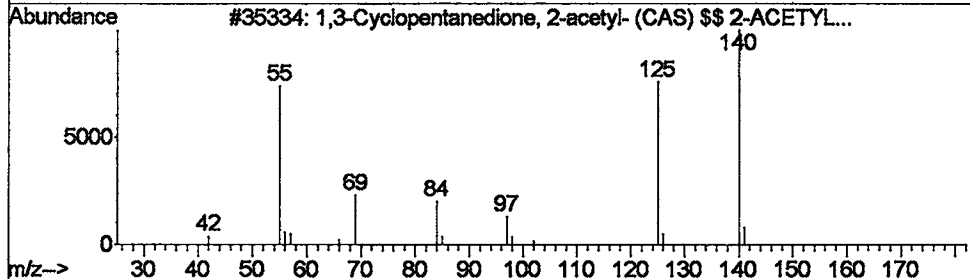
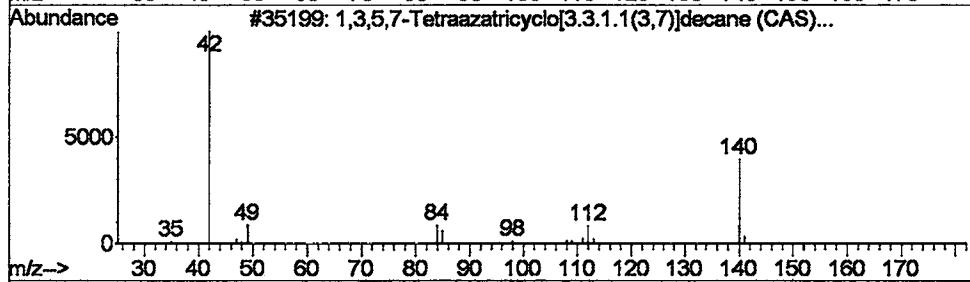
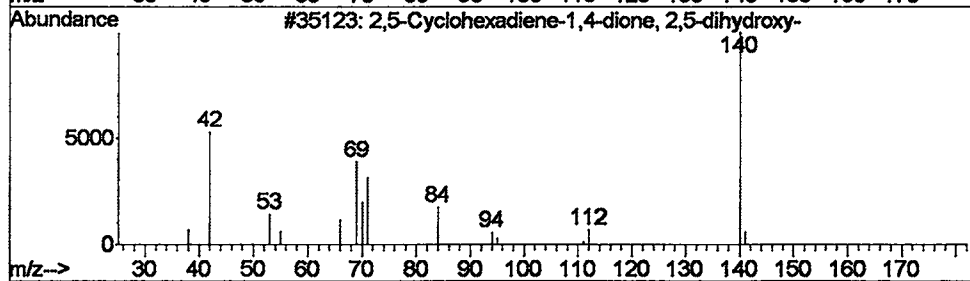
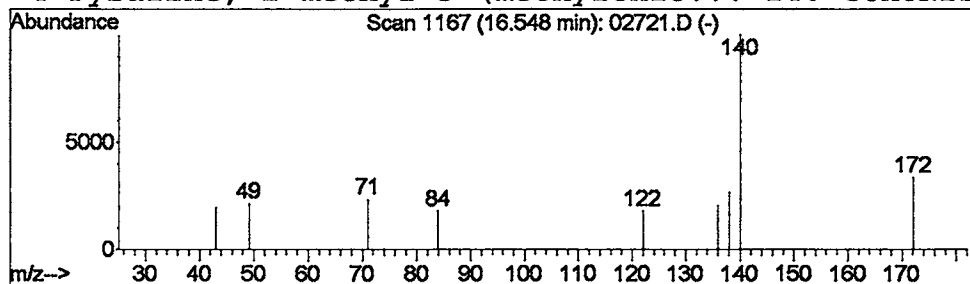
Title : 508 scan method

Library : C:\DATABASE\WILEY7N.L

Peak Number 8 2,5-Cyclohexadiene-1,4-dion... Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.55	0.03 ug/L	19813	Acenaphthene-d10	18.33

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2,5-Cyclohexadiene-1,4-dione, 2,...	140	C6H4O4	000615-94-1	9
2			1,3,5,7-Tetraazatricyclo[3.3.1.1...	140	C6H12N4	000100-97-0	7
3			1,3-Cyclopentanedione, 2-acetyl-...	140	C7H8O3	003859-39-0	5
4			Pyrazine, 2-methyl-3-(methylthio...	140	C6H8N2S	002882-20-4	5



Library Search Compound Report

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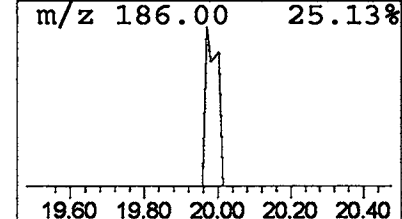
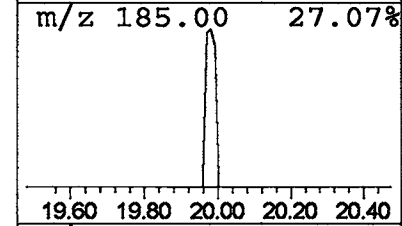
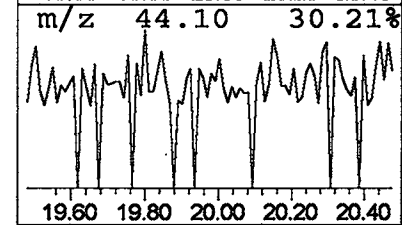
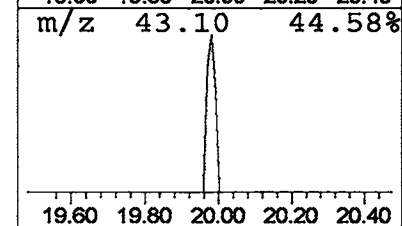
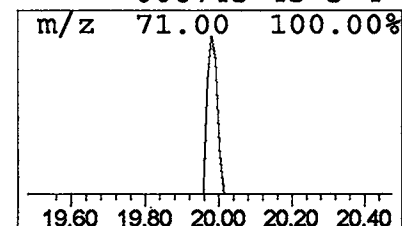
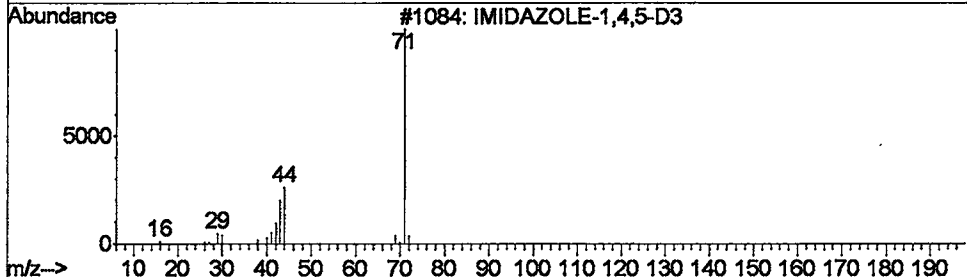
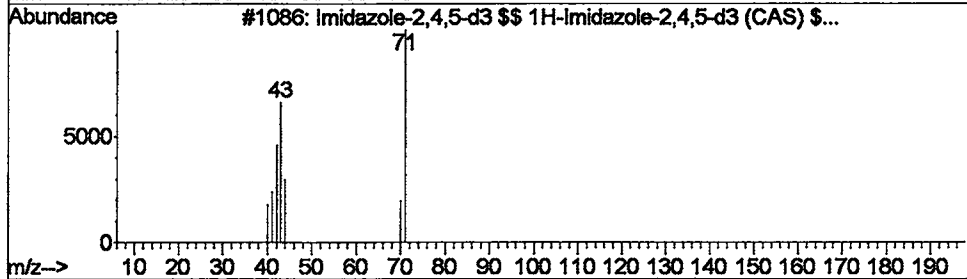
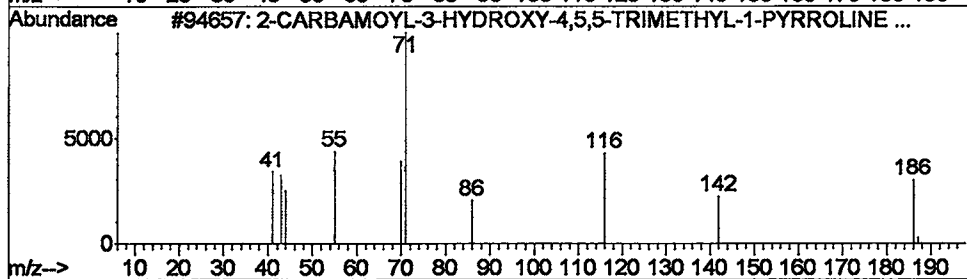
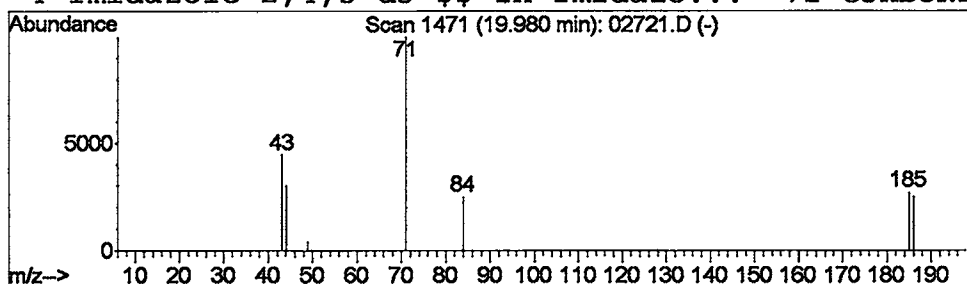
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 Inst : Instrumen
 Multiplr: 1.00

Quant Method : C:\MSDCHEM\1...\HP022102.M (RTE Integrator)
 Title : 508 scan method
 Library : C:\DATABASE\WILEY7N.L

 Peak Number 9 2-CARBAMOYL-3-HYDROXY-4,5,5... Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.98	0.03 ug/L	17003	Acenaphthene-d10	18.33

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-CARBAMOYL-3-HYDROXY-4,5,5-TRIM...	186	C8H14N2O3	072700-98-2	9
2			Imidazole-2,4,5-d3 \$\$ 1H-Imidazo...	71	C3HD3N2	006745-43-3	7
3			IMIDAZOLE-1,4,5-D3	71	C3HD3N2	053716-57-7	4
4			Imidazole-2,4,5-d3 \$\$ 1H-Imidazo...	71	C3HD3N2	006745-43-3	4



Library Search Compound Report

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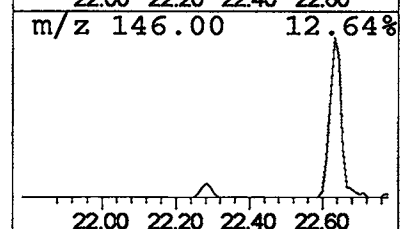
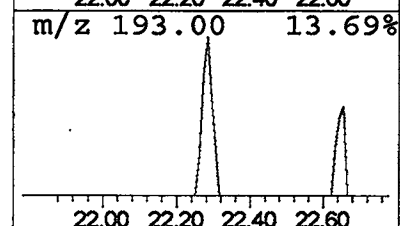
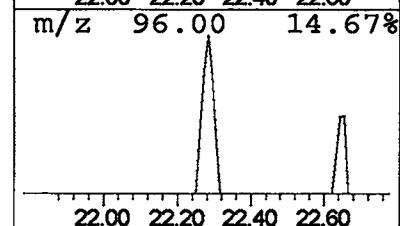
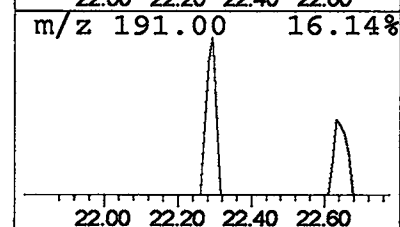
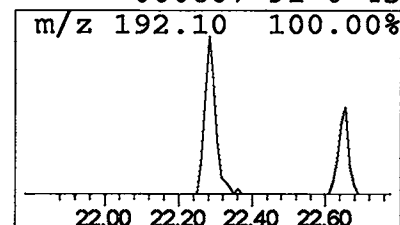
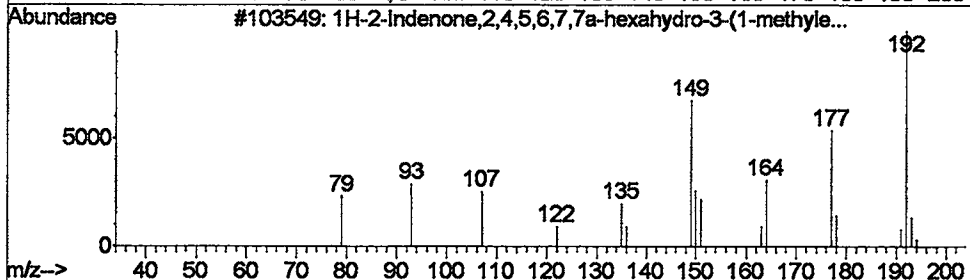
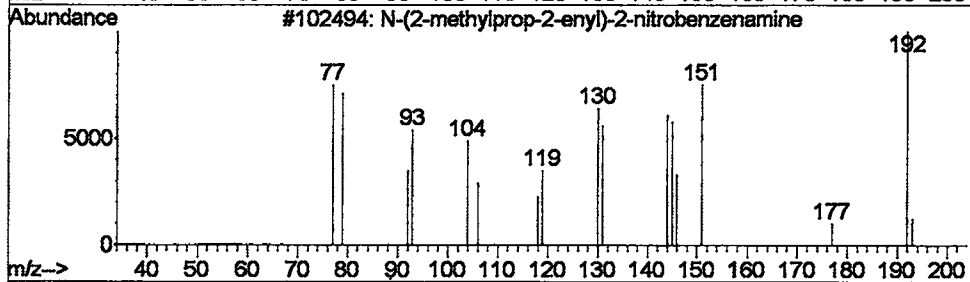
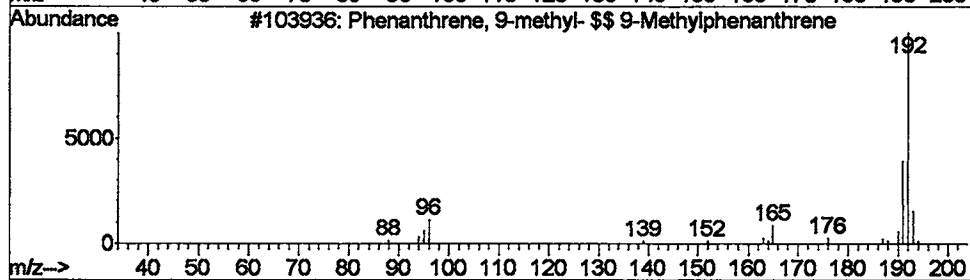
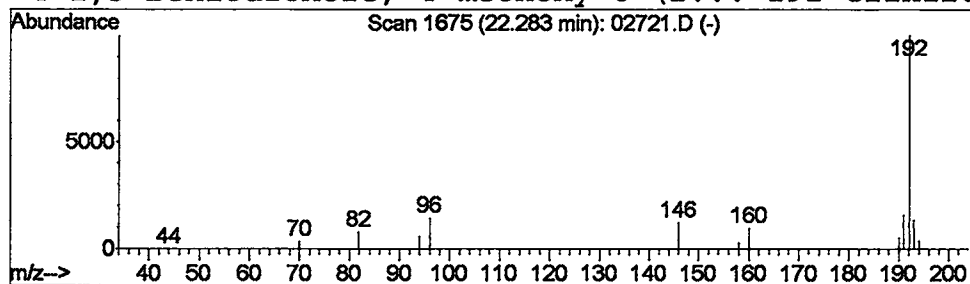
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Quant Method : C:\MSDCHEM\1...\HP022102.M (RTE Integrator)
 Title : 508 scan method
 Library : C:\DATABASE\WILEY7N.L

 Peak Number 10 Phenanthrene, 9-methyl- \$\$... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.28	0.11 ug/L	76233	Phenanthrene-d10	22.63

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Phenanthrene, 9-methyl- \$\$ 9-Met...	192	C15H12	000883-20-5	59
2			N-(2-methylprop-2-enyl)-2-nitrob...	192	C10H12N2O2	000000-00-0	52
3			1H-2-Indenone,2,4,5,6,7,7a-hexah...	192	C13H20O	000000-00-0	45
4			1,3-Benzodioxole, 4-methoxy-6-(2...	192	C11H12O3	000607-91-0	45



Library Search Compound Report

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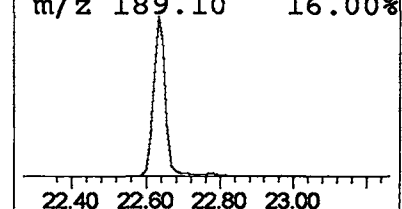
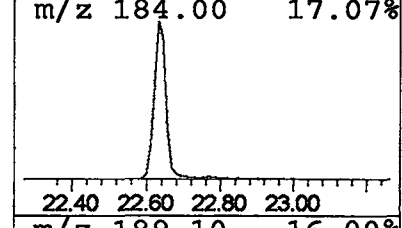
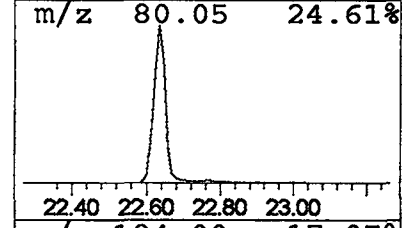
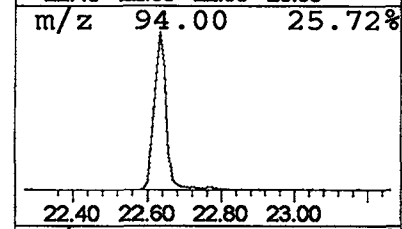
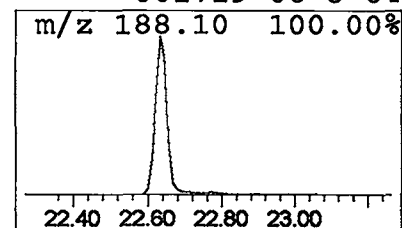
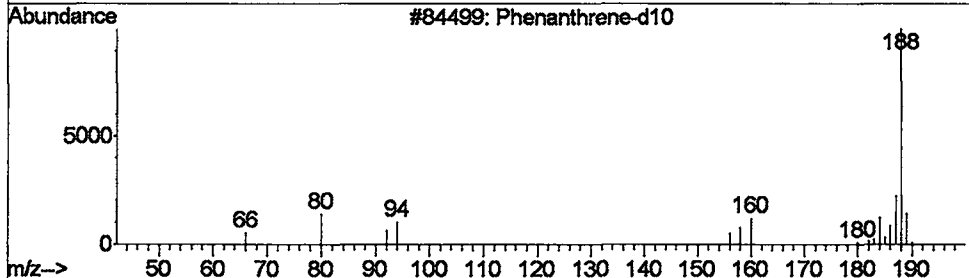
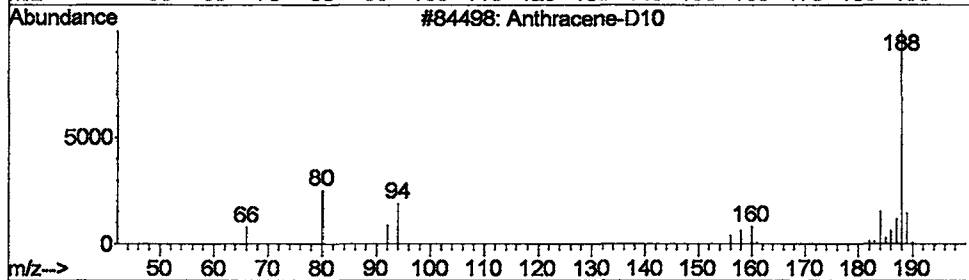
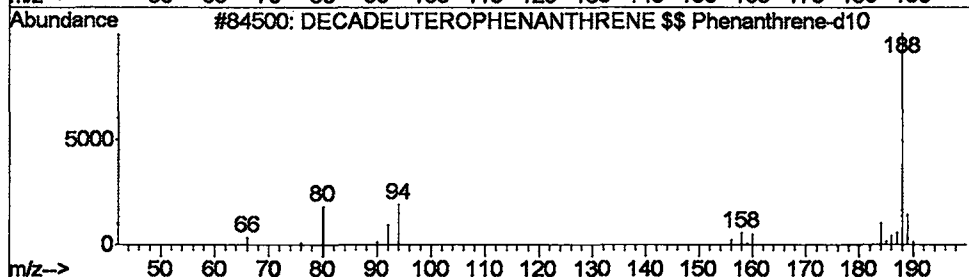
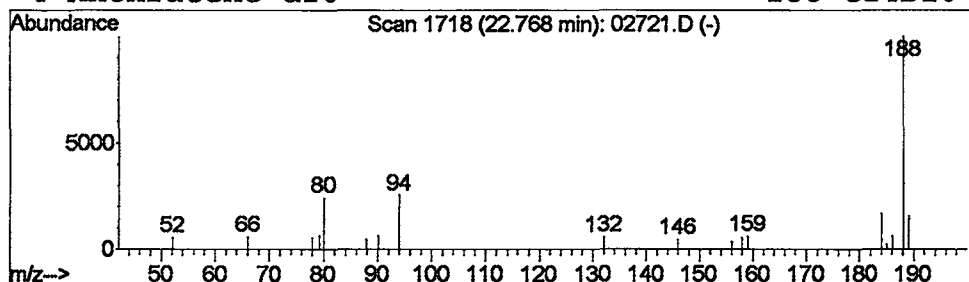
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 Inst : Instrumen
 Multiplr: 1.00

Quant Method : C:\MSDCHEM\1...\HP022102.M (RTE Integrator)
 Title : 508 scan method
 Library : C:\DATABASE\WILEY7N.L

 Peak Number 11 DECADEUTEROPHENANTHRENE \$\$... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.77	0.45 ug/L	301511	Phenanthrene-d10	22.63

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			DECADEUTEROPHENANTHRENE \$\$ Phena...	188	C14D10	001517-22-2	91
2			Anthracene-D10	188	C14D10	000000-00-0	87
3			Phenanthrene-d10	188	C14D10	001517-22-2	72
4			Anthracene-d10-	188	C14D10	001719-06-8	64



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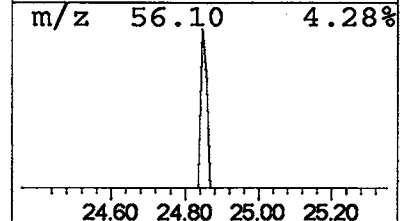
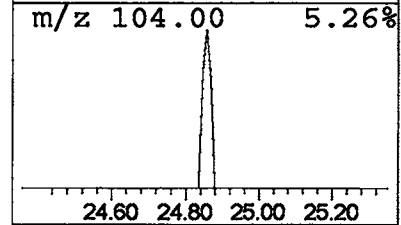
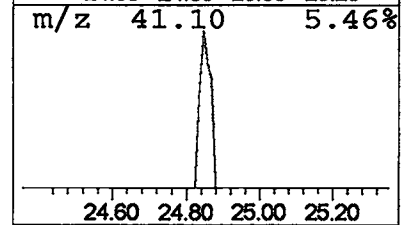
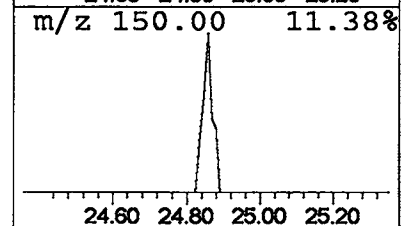
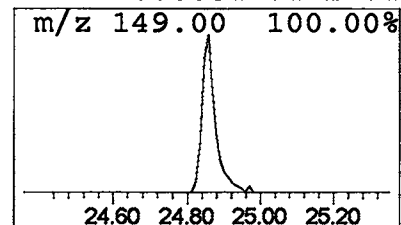
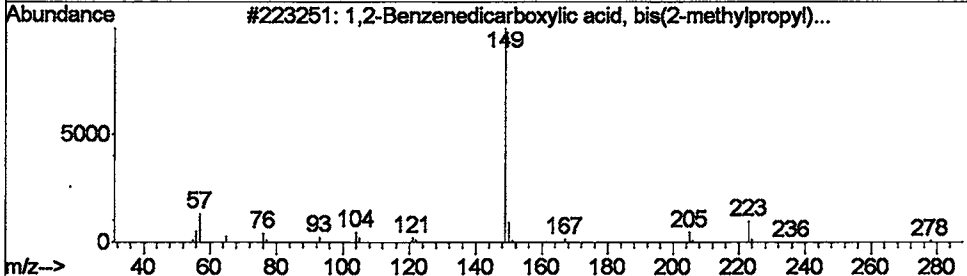
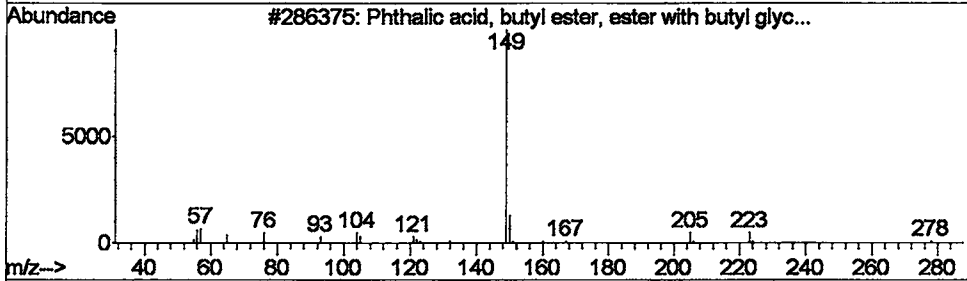
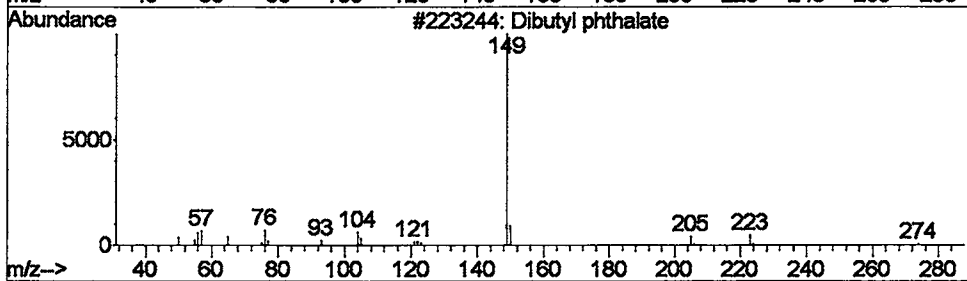
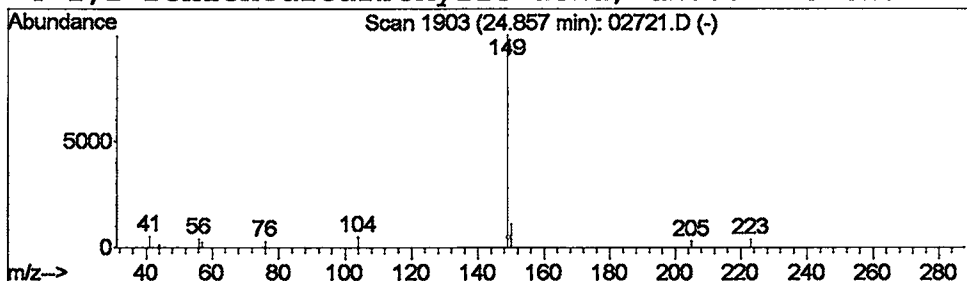
Vial: 6
 Operator:
 Inst : Instrumen
 Multiplr: 1.00

Quant Method : C:\MSDCHEM\1...\HP022102.M (RTE Integrator)
 Title : 508 scan method
 Library : C:\DATABASE\WILEY7N.L

 Peak Number 12 Dibutyl phthalate Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.86	0.10 ug/L	66389	Phenanthrene-d10	22.63

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Dibutyl phthalate	278	C16H22O4	000084-74-2	74
2			Phthalic acid, butyl ester, este...	336	C18H24O6	000085-70-1	74
3			1,2-Benzenedicarboxylic acid, bi...	278	C16H22O4	000084-69-5	74
4			1,2-Benzenedicarboxylic acid, di...	278	C16H22O4	000084-74-2	74



Library Search Compound Report

* Data File : C:\MSDCHEM\1\5973N\02FEB28\02721.D
Acq On : 28 Feb 2002 6:50 pm
Sample : 508 scan
Misc : February 28, 2002
MS Integration Params: LSCINT.P

Vial: 6
Operator:
Inst : Instrumen
Multiplr: 1.00

Quant Method : C:\MSDCHEM\1...\HP022102.M (RTE Integrator)

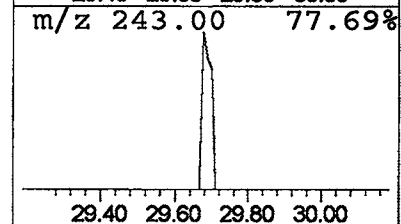
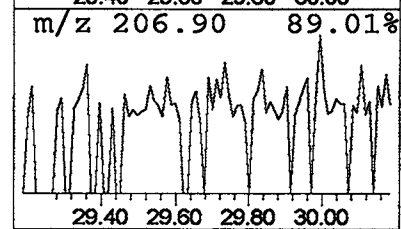
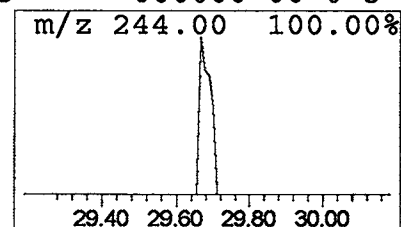
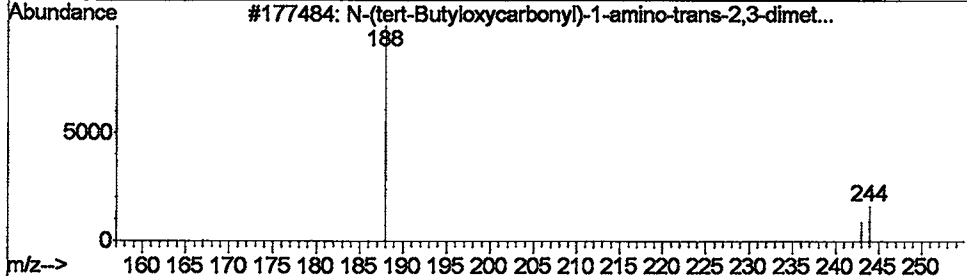
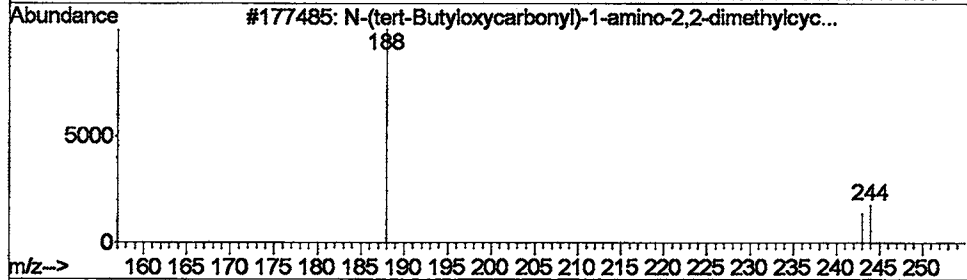
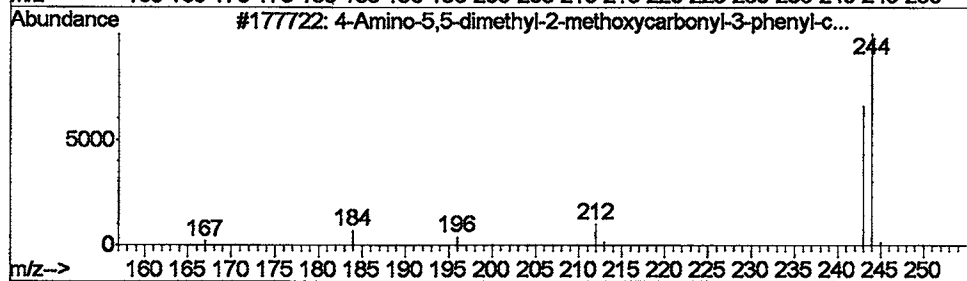
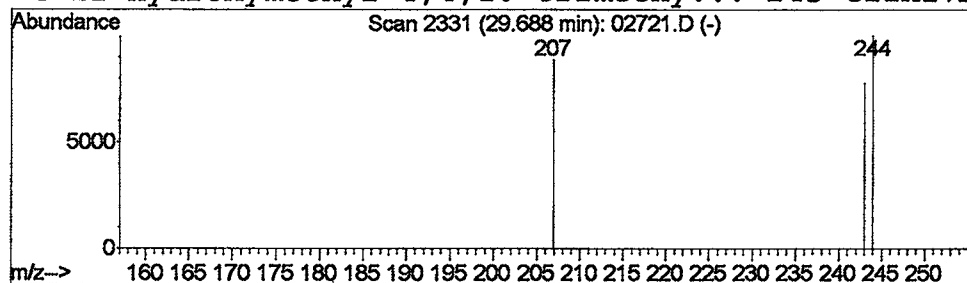
Title : 508 scan method

Library : C:\DATABASE\WILEY7N.L

Peak Number 13 4-Amino-5,5-dimethyl-2-meth... Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
29.69	0.03 ug/L	15726	Chrysene-d12	30.49

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		4-Amino-5,5-dimethyl-2-methoxyca...	243	C15H17NO2	118089-29-5	5
2		N-(tert-Butyloxycarbonyl)-1-amin...	243	C12H21NO4	000000-00-0	4
3		N-(tert-Butyloxycarbonyl)-1-amin...	243	C12H21NO4	000000-00-0	4
4		11-Hydroxymethyl-4,4,10-trimethy...	243	C11H17NO5	000000-00-0	3



Tentatively Identified Compound (LSC) summary

Operator ID: Date Acquired: 28 Feb 2002 6:50 pm
 Data File: C:\MSDCHEM\1\5973N\02FEB28\02721.D
 Name: 508 scan
 Misc: February 28, 2002
 Method: C:\MSDCHEM\1\METHODS\METHODS\HP022102.M (RTE Integrator)
 Title: 508 scan method
 Library Searched: C:\DATABASE\WILEY7N.L

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Butanoic acid, me...	3.62	0.1	ug/L	52411	1	9.72	7966840	20.0
ISOBUTYRALDEHYDE,...	3.93	0.1	ug/L	22817	1	9.72	7966840	20.0
Cyclobutene, 2-pr...	4.38	0.0	ug/L	18429	1	9.72	7966840	20.0
3-(CYCLOHEX-3'-EN...	5.94	0.3	ug/L	134636	1	9.72	7966840	20.0
4-Cyanocyclohexen...	9.96	0.0	ug/L	19370	1	9.72	7966840	20.0
Cyclopentasiloxan...	12.54	0.0	ug/L	26953	2	13.20	11618600	20.0
METHYL-2-DEOXY-L-...	13.39	0.1	ug/L	38017	2	13.20	11618600	20.0
2,5-Cyclohexadien...	16.55	0.0	ug/L	19813	3	18.33	12647900	20.0
2-CARBAMOYL-3-HYD...	19.98	0.0	ug/L	17003	3	18.33	12647900	20.0
Phenanthrene, 9-m...	22.28	0.1	ug/L	76233	4	22.63	13526100	20.0
DECADEUTEROPHENAN...	22.77	0.4	ug/L	301511	4	22.63	13526100	20.0
Dibutyl phthalate	24.86	0.1	ug/L	66389	4	22.63	13526100	20.0
4-Amino-5,5-dimet...	29.69	0.0	ug/L	15726	5	30.49	12059700	20.0