

User: Vinci, David L
Westchester Dept. of Labs & Research
Date: 03/15/2002 Time: 17:40:42

Sample: AE03184
Analysis: \$GCMS GCMS Scan For Unknowns
Units: ug
Started: 3/15/02 17:39 Ended: 3/15/02 17:39 Analyst: DLV
Page: 1

	Component Name	Viol	Result	Qualifier	MDL
1	Other unknown compounds		.		
2	Surr_2,4-DDD		117		5.0

Comments for Sample AE03184 - Analysis \$GCMS

Westchester Dept. of Labs & Research

User: Vinci, David L

Date: 03-15-2002 Time: 17:41:24

The GCMS scan, from 35 to 550 amu, does not reveal the presence of unusual compounds. The recoveries for 5 of the 43 compounds in the LCS Standard were outside their acceptance ranges.

Data File : C:\MSDCHEM\1\5973N\02MAR01\03184.D

Vial: 4

Acq On : 1 Mar 2002 12:48 pm

Operator:

Sample : 508 scan

Inst : Instrumen

Misc : March 1, 2002

Multiplr: 1.00

MS Integration Params: SPLIT.P

Quant Time: Mar 04 10:50:44 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	9.71	152	1264309	20.00	ug/L	0.00
10) Naphthalene-d8	13.20	136	5440371	20.00	ug/L	0.00
17) Acenaphthene-d10	18.33	164	2756311	20.00	ug/L	0.00
29) Phenanthrene-d10	22.63	188	4687831	20.00	ug/L	-0.01
38) Chrysene-d12	30.49	240	3802666	20.00	ug/L	-0.03
44) Perylene-d12	34.41	264	2492357	20.00	ug/L	-0.04

System Monitoring Compounds

37) 2,4-DDD	27.72	235	364141	5.83	ug/L	0.00
Spiked Amount	5.000		Recovery	=	116.60%	

Target Compounds

						Qvalue
20) Dimethylphthalate	18.33	163	444561	2.66	ug/L	# 59
21) 2,6-Dinitrotoluene	18.33	165	353777	9.16	ug/L	# 27

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

03184.D HP022102.M

Mon Mar 04 10:50:45 2002

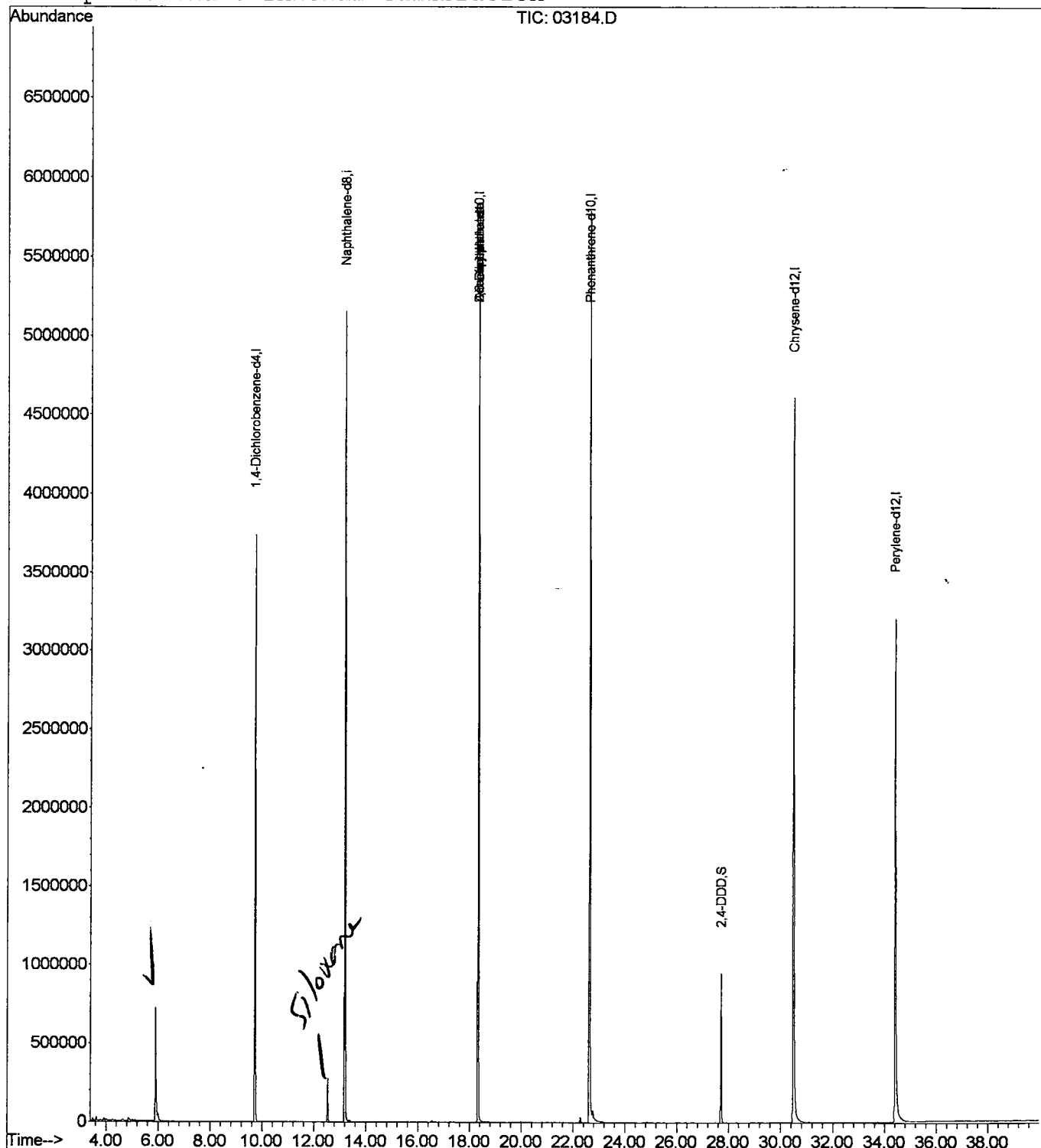
Page 1

Data File : C:\MSDCHEM\1\5973N\02MAR01\03184.D
 Acq On : 1 Mar 2002 12:48 pm
 Sample : 508 scan
 Misc : March 1, 2002
 MS Integration Params: SPLIT.P
 Quant Time: Mar 4 10:50 2002

Vial: 4
 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\02MAR01\03184.D

Vial: 4

Acq On : 1 Mar 2002 12:48 pm

Operator:

Sample : 508 scan

Inst : Instrumen

Misc : March 1, 2002

Multiplr: 1.00

MS Integration Params: LSCINT.P

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

Title : 508 scan method

Smoothing : OFF

Filtering: 5

Sampling : 2

Min Area: 100 Area counts

Start Thrs: 0.2

Max Peaks: 20

Stop Thrs : 0

Peak Location: TOP

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

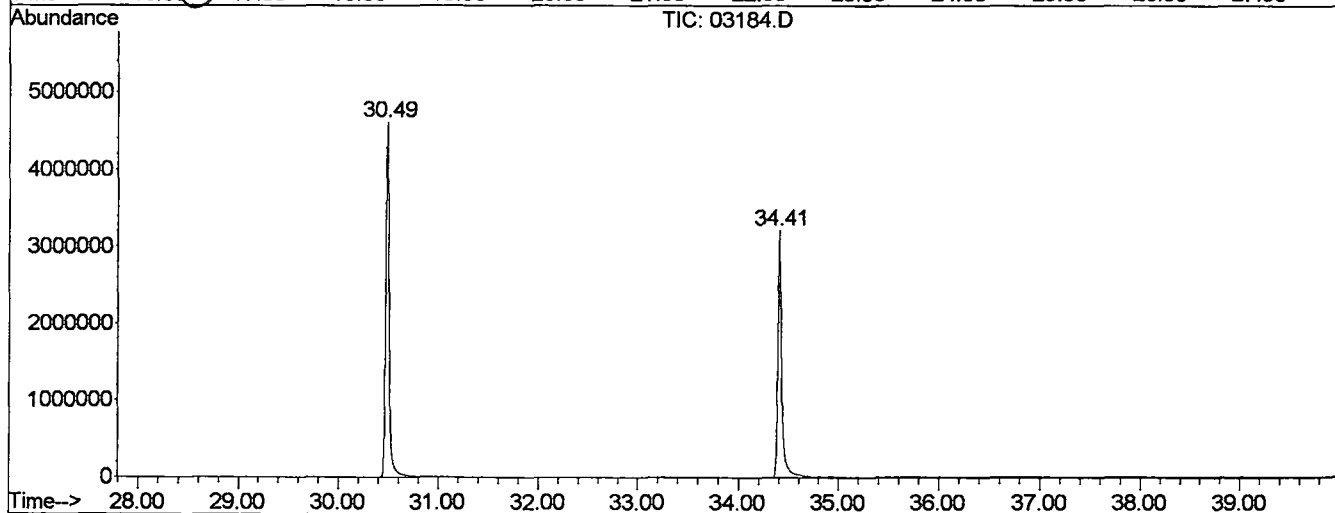
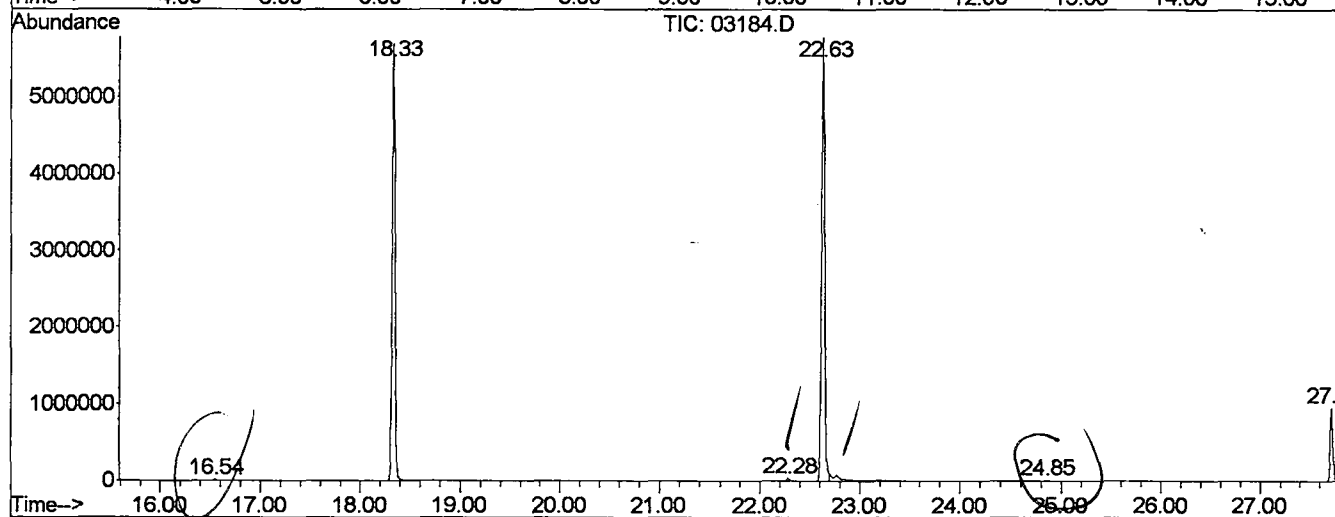
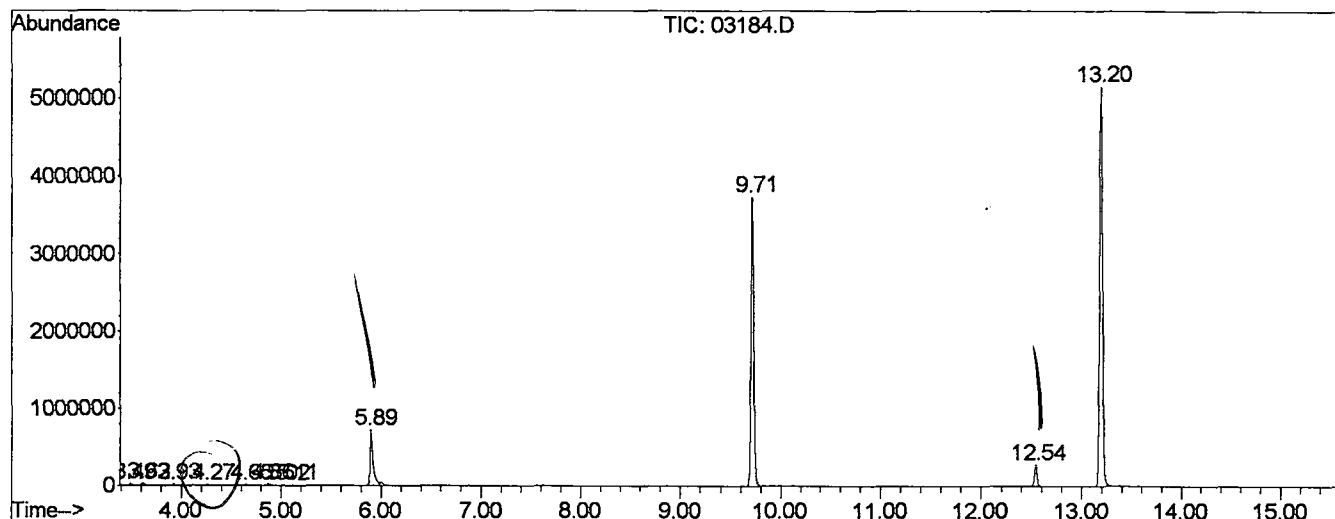
Signal : TIC

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.487	7	10	15	rVB	15513	24096	0.20%	0.037%
2	3.622	19	22	29	rBV2	24958	51419	0.42%	0.080%
3	3.927	45	49	53	rVV	15483	24714	0.20%	0.038%
4	4.266	75	79	85	rVB3	8957	21157	0.17%	0.033%
5	4.650	107	113	121	rBV5	10256	38457	0.32%	0.059%
6	4.864	127	132	143	rBV4	19619	81939	0.67%	0.127%
7	5.022	143	146	149	rVV2	10603	21857	0.18%	0.034%
8	5.112	151	154	165	rVV4	7776	20556	0.17%	0.032%
9	5.891	219	223	231	rBV	721114	1462430	12.00%	2.262%
10	9.707	557	561	567	rBV	3739008	7077223	58.08%	10.945%
11	12.541	807	812	817	rVB	277583	472050	3.87%	0.730%
12	13.195	865	870	875	rBV	5150577	10269013	84.27%	15.882%
13	16.537	1161	1166	1169	rBV	12573	25028	0.21%	0.039%
14	18.332	1319	1325	1331	rBV	5691749	11215318	92.03%	17.345%
15	22.283	1671	1675	1683	rBV	35522	72108	0.59%	0.112%
16	22.633	1701	1706	1711	rBV2	5785150	12186267	100.00%	18.847%
17	24.845	1899	1902	1911	rVB	8344	19955	0.16%	0.031%
18	27.724	2151	2157	2161	rBV	951393	1766673	14.50%	2.732%
19	30.490	2397	2402	2443	rVB	4605180	11233371	92.18%	17.373%
20	34.407	2743	2749	2791	rBV	3208231	8576550	70.38%	13.264%

Sum of corrected areas: 64660181

LSC Report - Integrated Chromatogram

File : C:\MSDCHEM\1\5973N\02MAR01\03184.D
 Operator :
 Acquired : 1 Mar 2002 12:48 pm using AcqMethod HP020702
 Instrument : Instrumen
 Sample Name: 508 scan
 Misc Info : March 1, 2002
 Vial Number: 4
 Quant File :HP022102.RES (RTE Integrator)



Library Search Compound Report

Data File : C:\MSDCHEM\1\5973N\02MAR01\03184.D
Acq On : 1 Mar 2002 12:48 pm
Sample : 508 scan
Misc : March 1, 2002
MS Integration Params: LSCINT.P

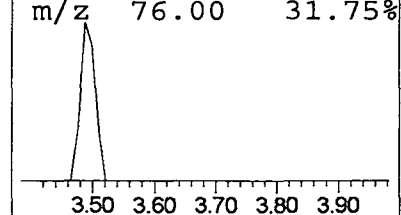
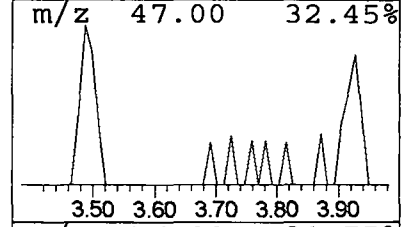
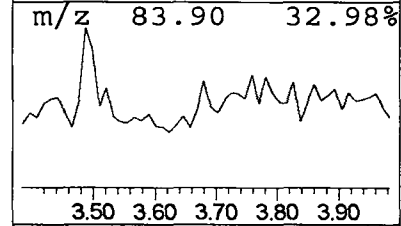
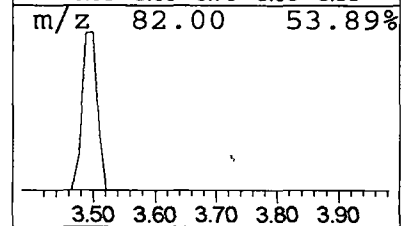
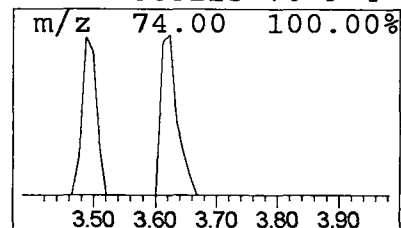
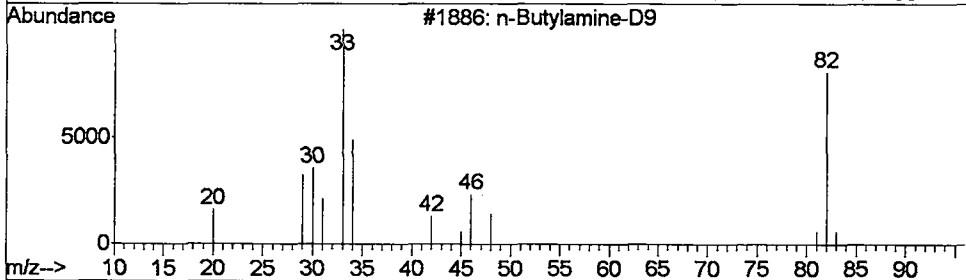
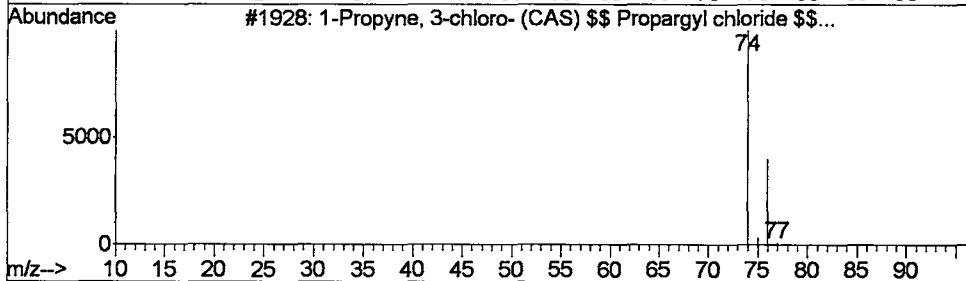
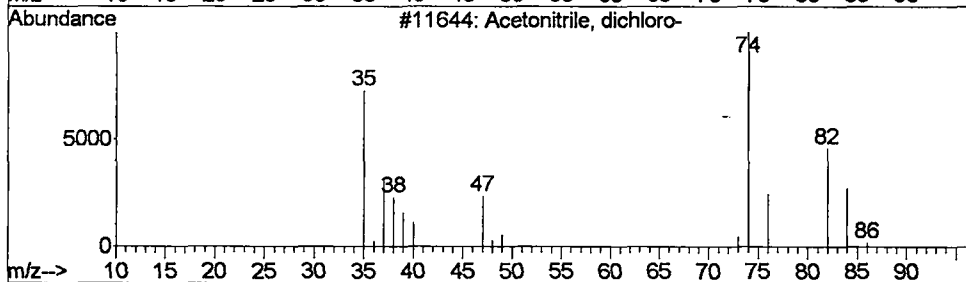
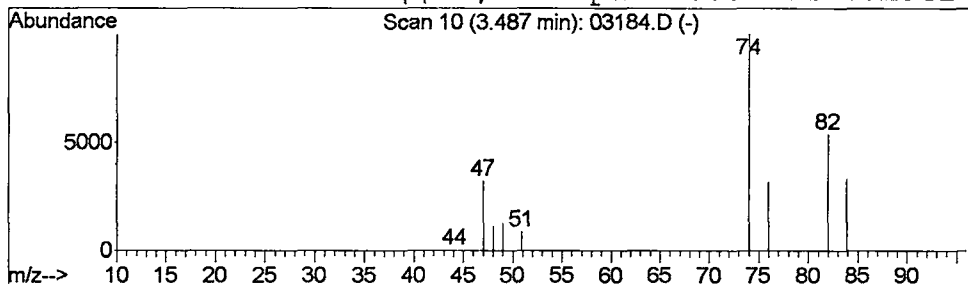
Vial: 4
Operator:
Inst : Instrumen
Multiplr: 1.00

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

Peak Number 1 Acetonitrile, dichloro- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.49	0.07 ug/L	24096	1,4-Dichlorobenzene-d4	9.71

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Acetonitrile, dichloro-	109	C2HCl2N	003018-12-0	83
2			1-Propyne, 3-chloro- (CAS) \$\$ Pr...	74	C3H3Cl	000624-65-7	5
3			n-Butylamine-D9	82	C4H2D9N	000000-00-0	5
4			CHLOROPROPADIENE \$\$ 1,2-Propadie...	74	C3H3Cl	003223-70-9	4



Library Search Compound Report

Data File : C:\MSDCHEM\1\5973N\02MAR01\03184.D
Acq On : 1 Mar 2002 12:48 pm
Sample : 508 scan
Misc : March 1, 2002
MS Integration Params: LSCINT.P

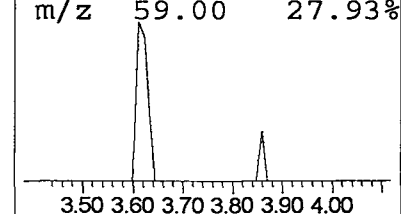
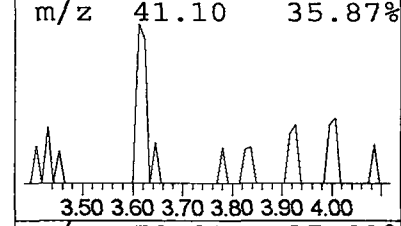
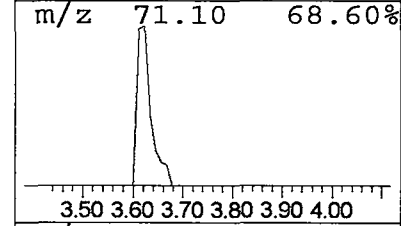
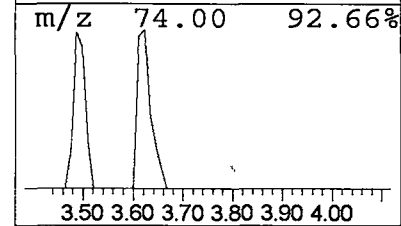
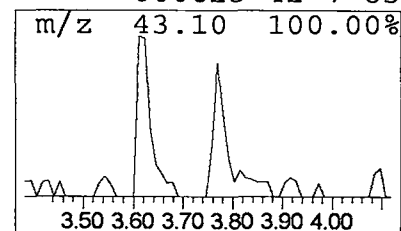
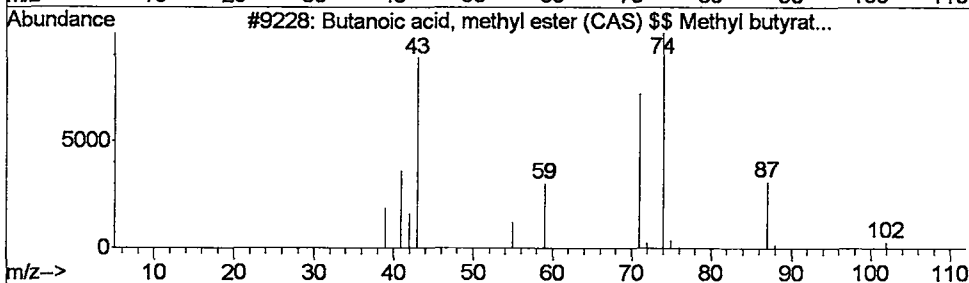
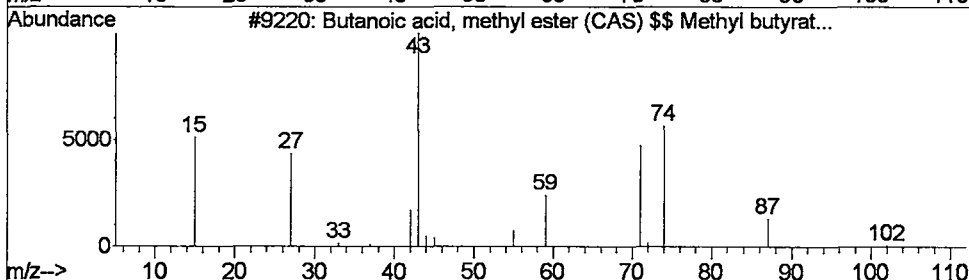
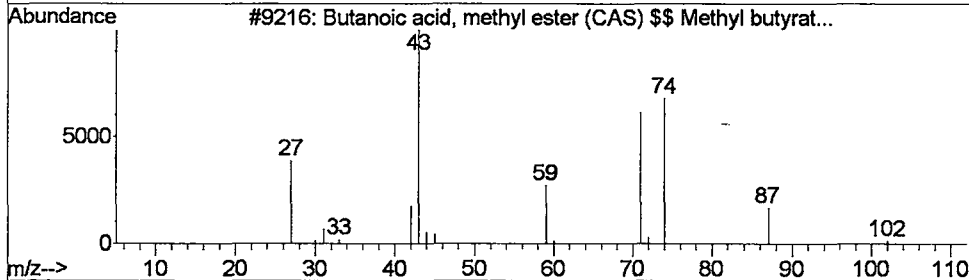
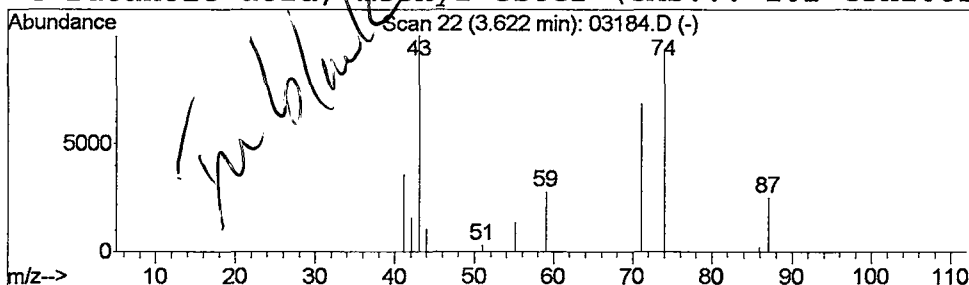
Vial: 4
Operator:
Inst : Instrumen
Multiplr: 1.00

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

Peak Number 2 Butanoic acid, methyl ester... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.62	0.15 ug/L	51419	1,4-Dichlorobenzene-d4	9.71

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



Library Search Compound Report

Data File : C:\MSDCHEM\1\5973N\02MAR01\03184.D
Acq On : 1 Mar 2002 12:48 pm
Sample : 508 scan
Misc : March 1, 2002
MS Integration Params: LSCINT.P

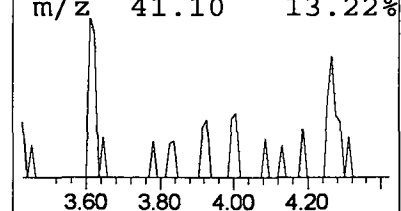
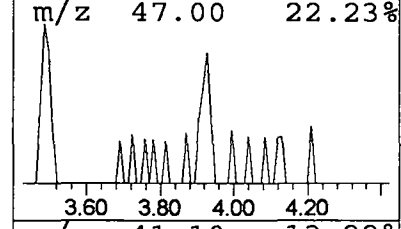
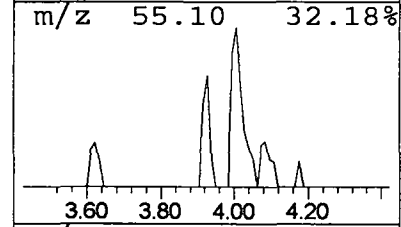
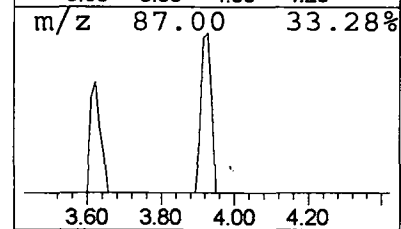
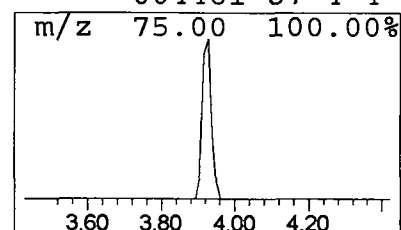
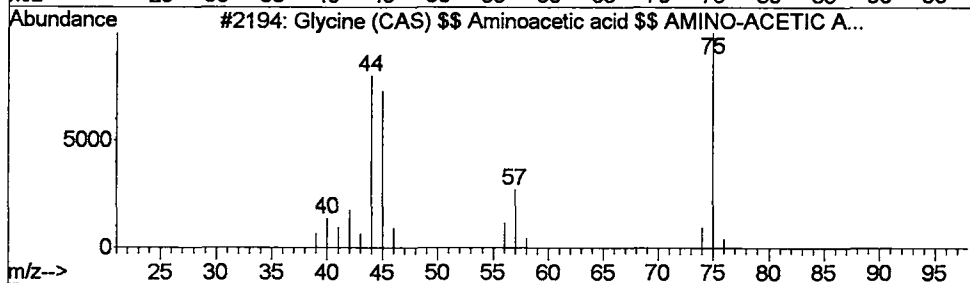
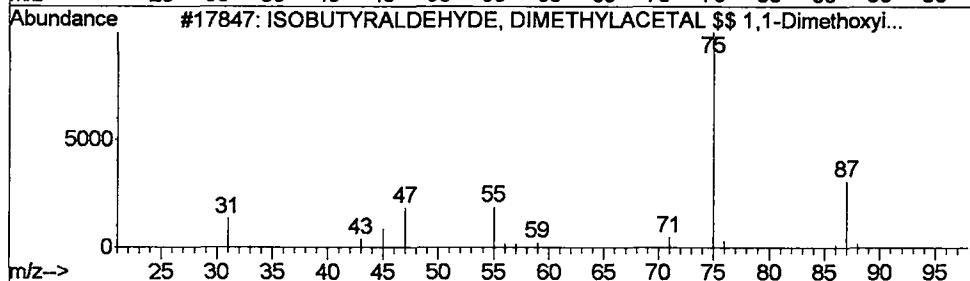
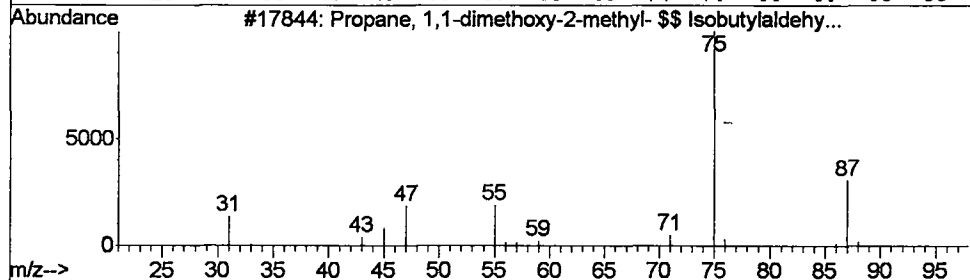
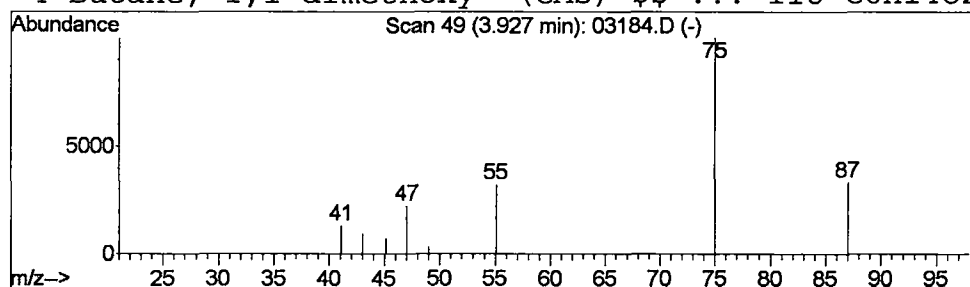
Vial: 4
Operator:
Inst : Instrumen
Multiplr: 1.00

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

Peak Number 3 Propane, 1,1-dimethoxy-2-me... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.93	0.07 ug/L	24714	1,4-Dichlorobenzene-d4	9.71

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Propane, 1,1-dimethoxy-2-methyl-...	118	C6H14O2	041632-89-7	9
2		ISOBUTYRALDEHYDE, DIMETHYLACETAL...	118	C6H14O2	000000-00-0	9
3		Glycine (CAS) \$\$ Aminoacetic aci...	75	C2H5NO2	000056-40-6	4
4		Butane, 1,1-dimethoxy- (CAS) \$\$...	118	C6H14O2	004461-87-4	4



Library Search Compound Report

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Acq On : 1 Mar 2002 12:48 pm
Sample : 508 scan
Misc : March 1, 2002
MS Integration Params: LSCINT.P

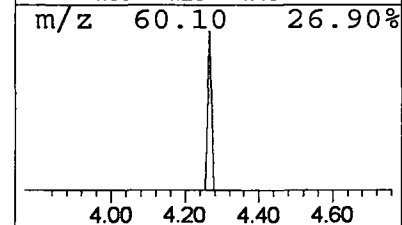
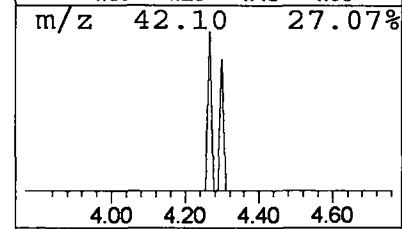
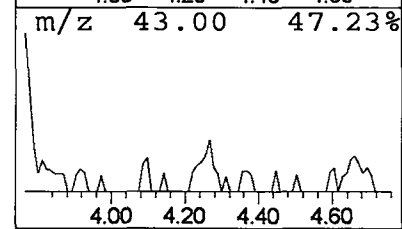
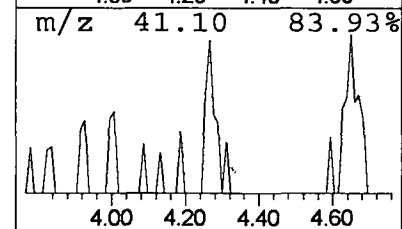
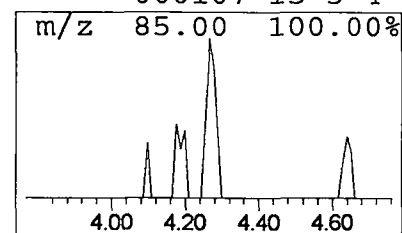
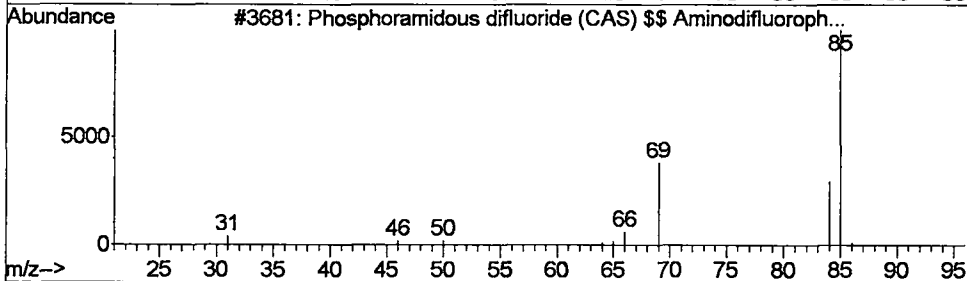
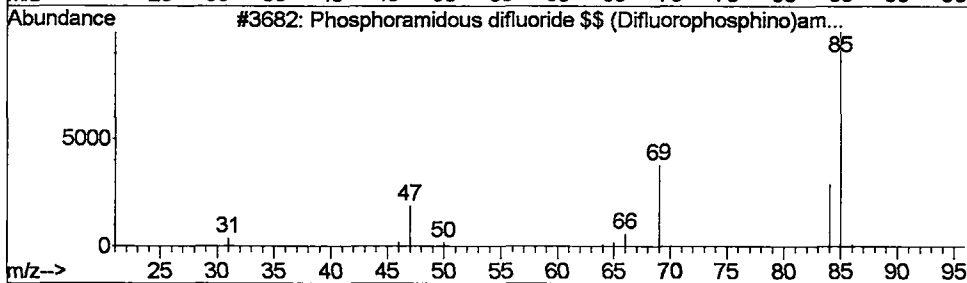
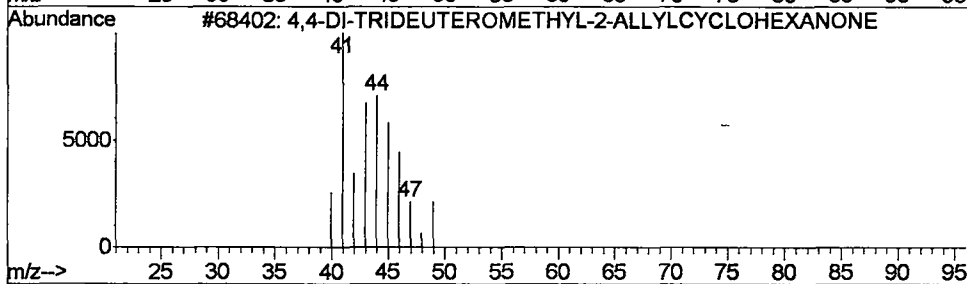
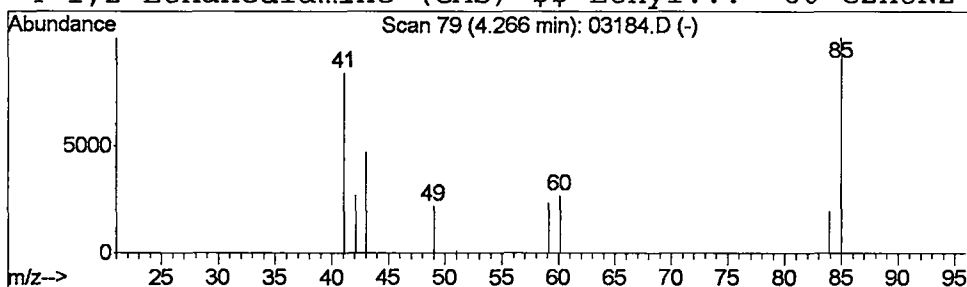
Vial: 4
Operator:
Inst : Instrumen
Multiplr: 1.00

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

Peak Number 4 4,4-DI-TRIDEUTEROMETHYL-2-A... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.27	0.06 ug/L	21157	1,4-Dichlorobenzene-d4	9.71

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			4,4-DI-TRIDEUTEROMETHYL-2-ALLYLC...	172	C11H12D6O	000000-00-0	9
2			Phosphoramidous difluoride \$\$ (D...	85	F2H2NP	025757-74-8	5
3			Phosphoramidous difluoride (CAS)...	85	F2H2NP	025757-74-8	5
4			1,2-Ethanediamine (CAS) \$\$ Ethyl...	60	C2H8N2	000107-15-3	4



Library Search Compound Report

Data File : C:\MSDCHEM\1\5973N\02MAR01\03184.D

Acq On : 1 Mar 2002 12:48 pm

Sample : 508 scan

Misc : March 1, 2002

MS Integration Params: LSCINT.P

Vial: 4

Operator:

Inst : Instrumen

Multiplr: 1.00

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

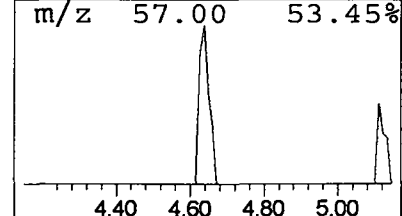
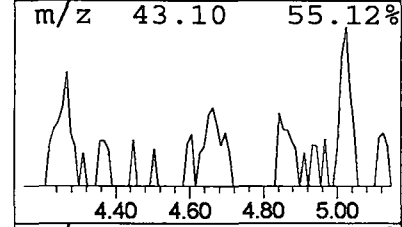
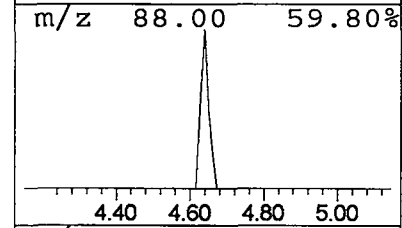
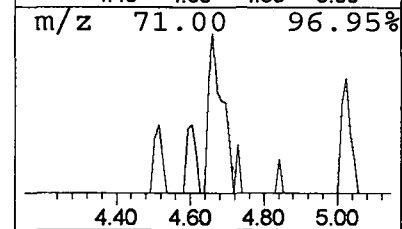
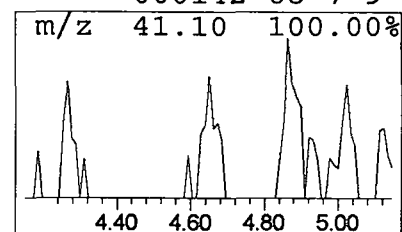
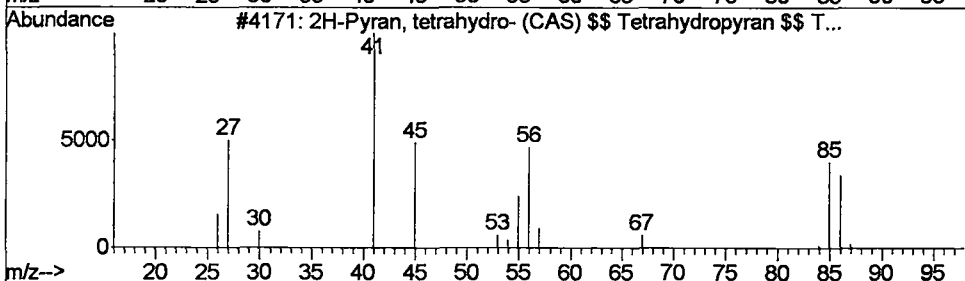
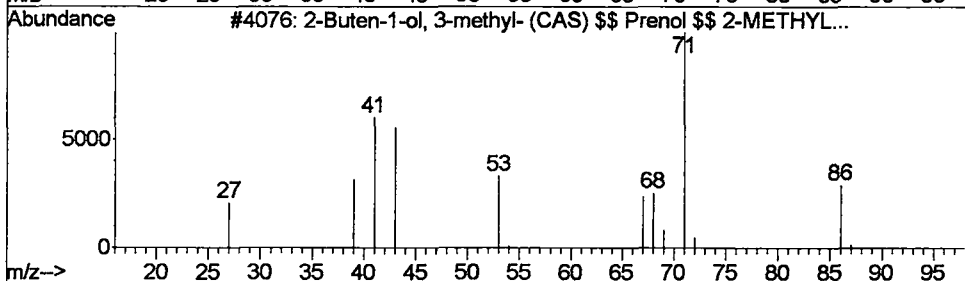
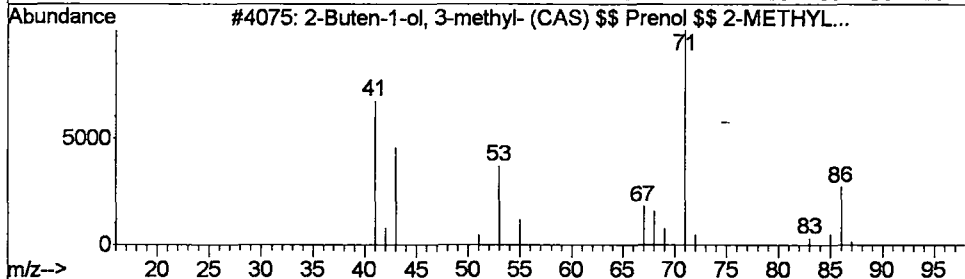
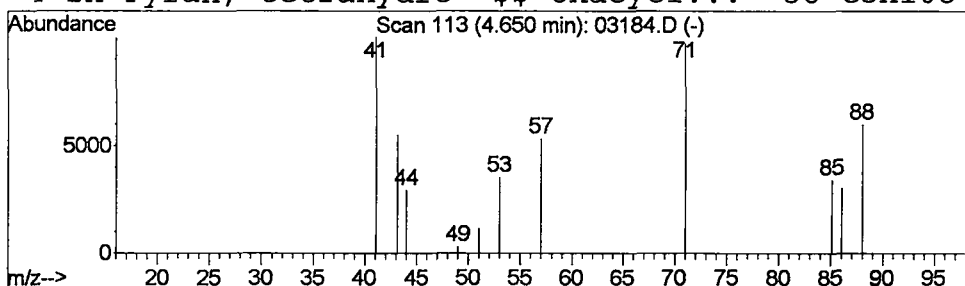
Title : 508 scan method

Library : C:\DATABASE\WILEY7N.L

Peak Number 5 2-Buten-1-ol, 3-methyl- (CA... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.65	0.11 ug/L	38457	1,4-Dichlorobenzene-d4	9.71

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Buten-1-ol, 3-methyl- (CAS) \$\$. . .	86	C5H10O	000556-82-1	47
2			2-Buten-1-ol, 3-methyl- (CAS) \$\$. . .	86	C5H10O	000556-82-1	37
3			2H-Pyran, tetrahydro- (CAS) \$\$. T. . .	86	C5H10O	000142-68-7	9
4			2H-Pyran, tetrahydro- \$\$. Oxacycl. . .	86	C5H10O	000142-68-7	9



Library Search Compound Report

Data File : C:\MSDCHEM\1\5973N\02MAR01\03184.D
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Misc : March 1, 2002
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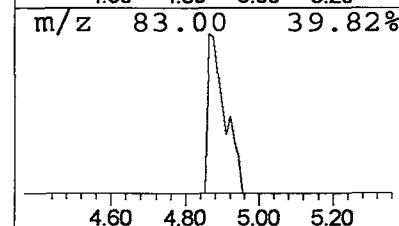
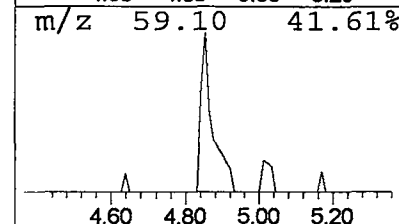
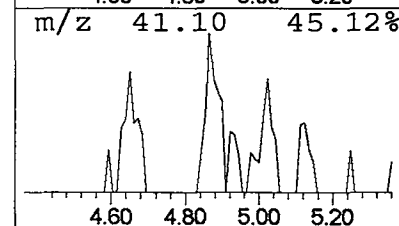
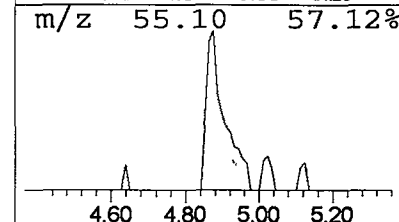
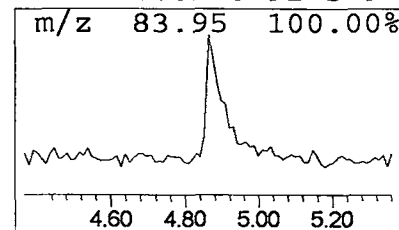
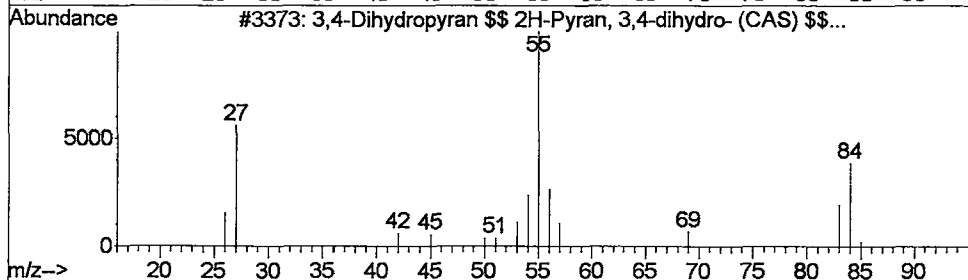
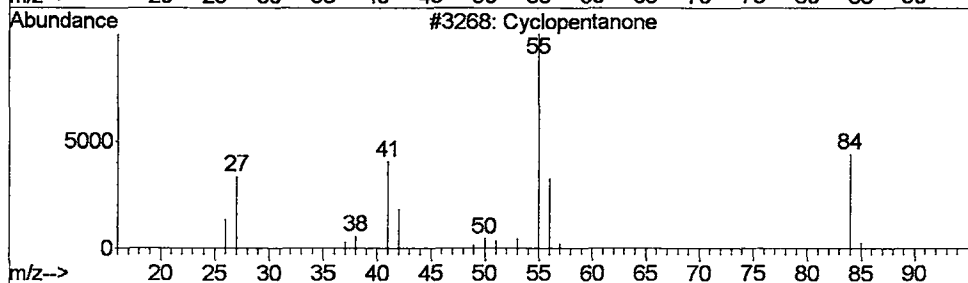
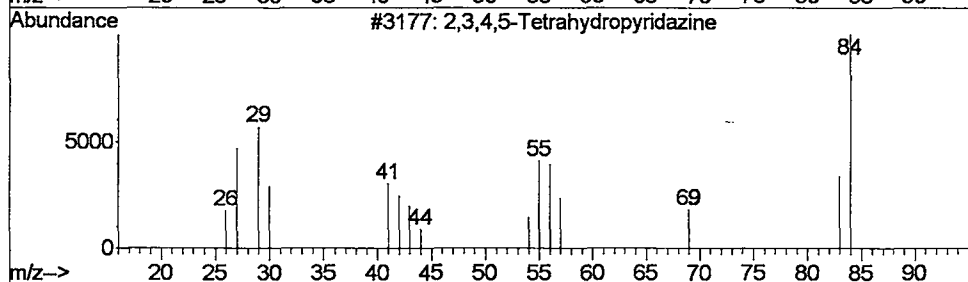
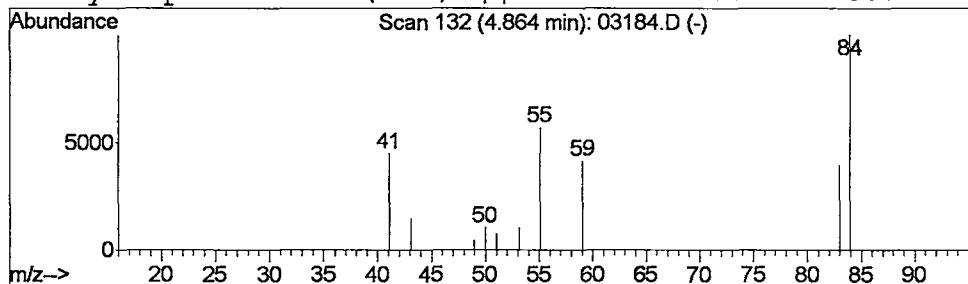
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Quant Method : C:\MSDCHEM\1...\HP022102.M (RTE Integrator)
Title : 508 scan method
Library : C:\DATABASE\WILEY7N.L

Peak Number 6 2,3,4,5-Tetrahydropyridazine Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.86	0.23 ug/L	81939	1,4-Dichlorobenzene-d4	9.71

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2,3,4,5-Tetrahydropyridazine	84	C4H8N2	000694-06-4	42
2			Cyclopentanone	84	C5H8O	000120-92-3	32
3			3,4-Dihydropyran \$\$ 2H-Pyran, 3,...	84	C5H8O	000110-87-2	9
4			Cyclopentanone (CAS) \$\$ Dumasin ...	84	C5H8O	000120-92-3	9



Library Search Compound Report

Data File : C:\MSDCHEM\1\5973N\02MAR01\03184.D
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Misc : March 1, 2002
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Multiplr: 1.00

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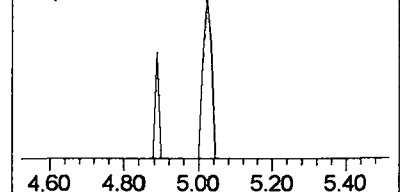
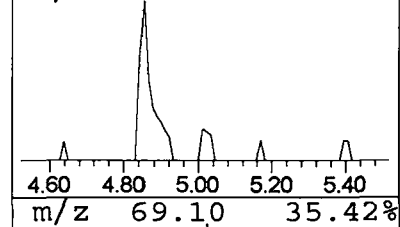
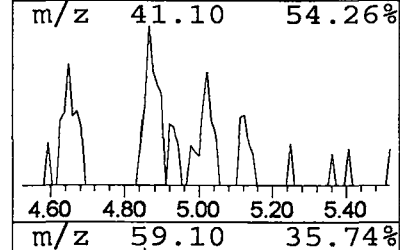
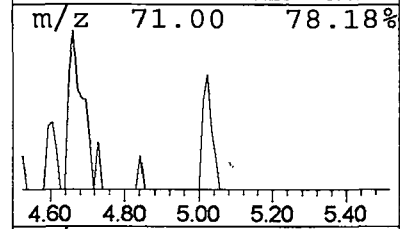
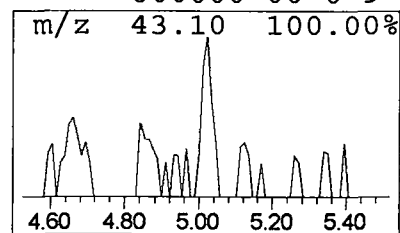
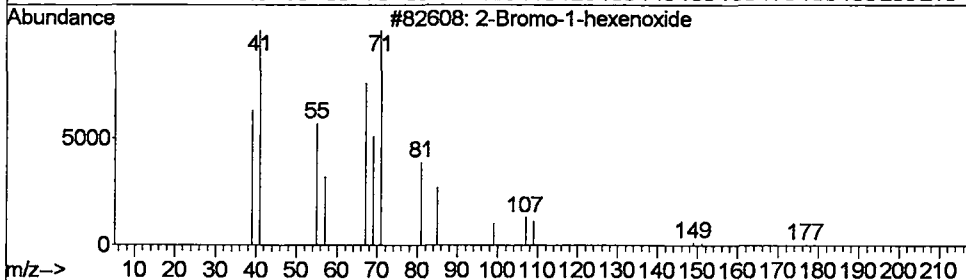
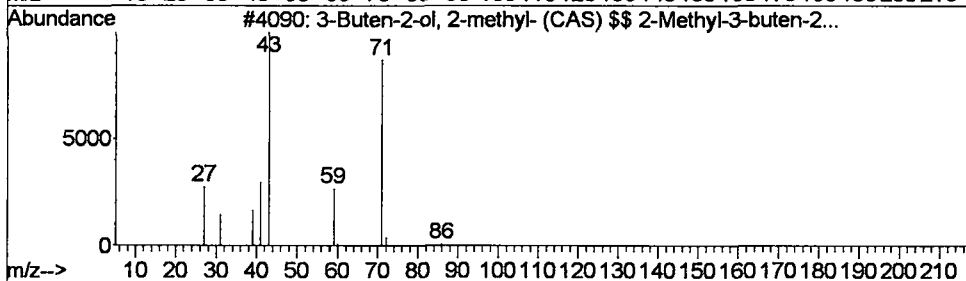
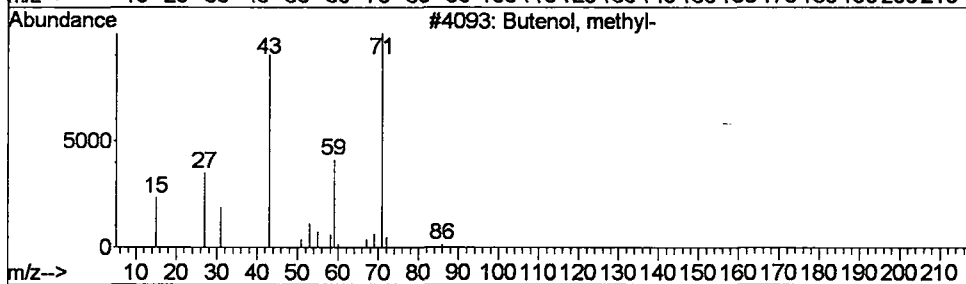
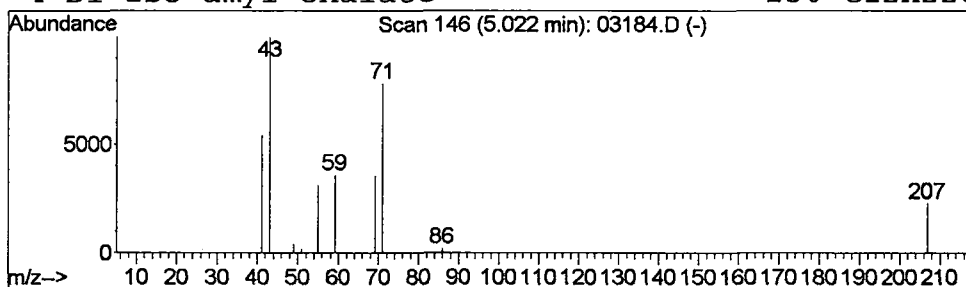
Title : 508 scan method

Library : C:\DATABASE\WILEY7N.L

Peak Number 7 Butenol, methyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.02	0.06 ug/L	21857	1,4-Dichlorobenzene-d4	9.71

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Butenol, methyl-	86	C5H10O	060766-00-9	50
2		3-Buten-2-ol, 2-methyl- (CAS) \$\$...	86	C5H10O	000115-18-4	9
3		2-Bromo-1-hexenoxide	178	C6H11BrO	000000-00-0	9
4		Di-iso-amyl oxalate	230	C12H22O4	000000-00-0	9



Library Search Compound Report

Data File : C:\MSDCHEM\1\5973N\02MAR01\03184.D

Acq On : 1 Mar 2002 12:48 pm

Sample : 508 scan

Misc : March 1, 2002

MS Integration Params: LSCINT.P

Vial: 4

Operator:

Inst : Instrumen

Multiplr: 1.00

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

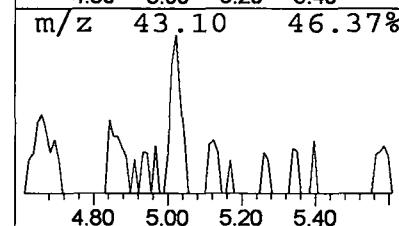
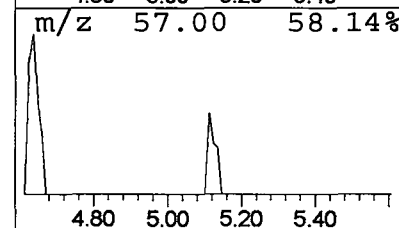
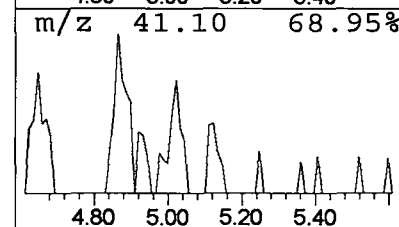
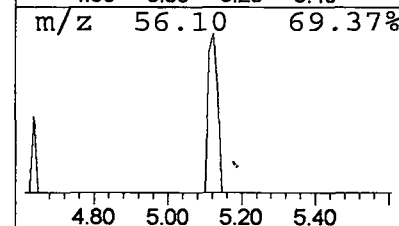
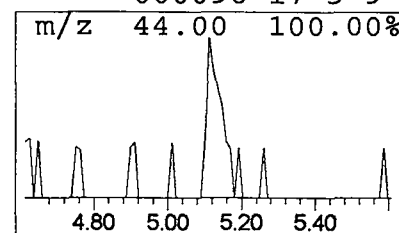
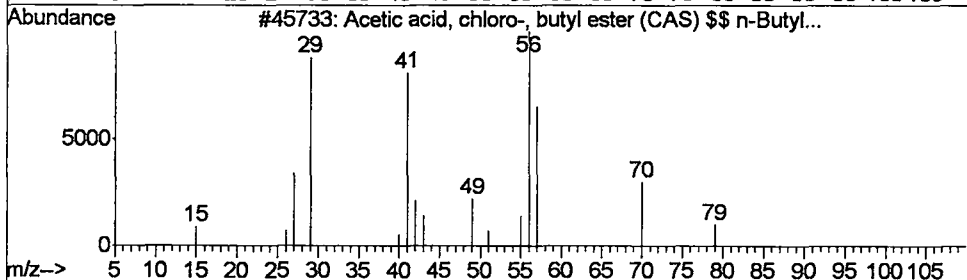
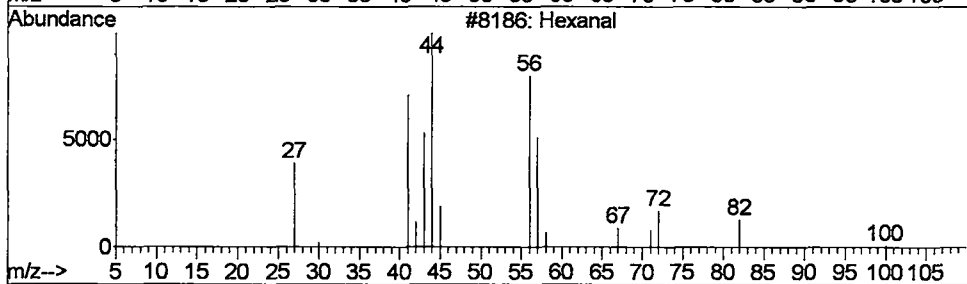
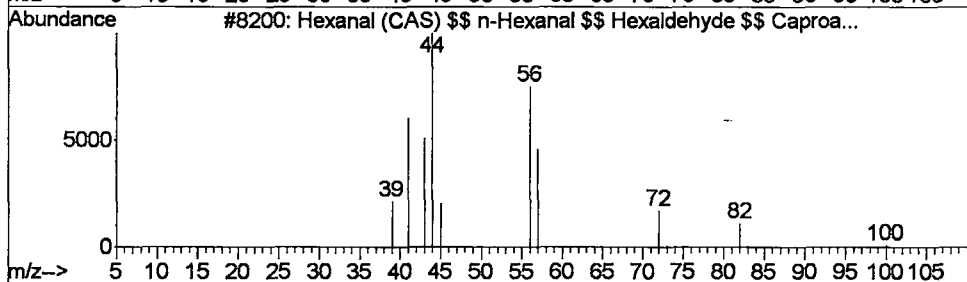
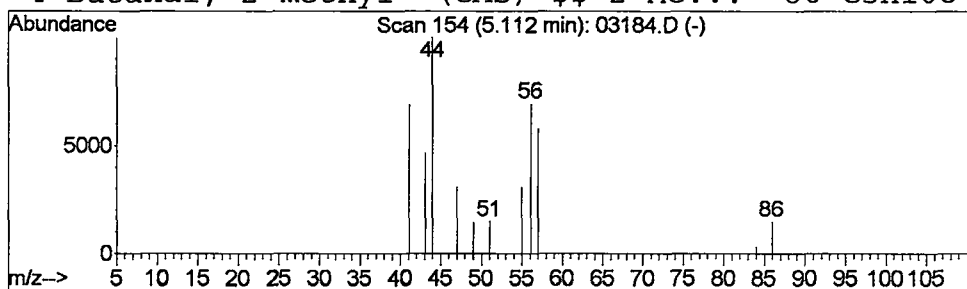
Title : 508 scan method

Library : C:\DATABASE\WILEY7N.L

Peak Number 8 Hexanal (CAS) \$\$ n-Hexanal ... Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.11	0.06 ug/L	20556	1,4-Dichlorobenzene-d4	9.71

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Hexanal (CAS) \$\$ n-Hexanal \$\$ He...	100	C6H12O	000066-25-1	38
2		Hexanal	100	C6H12O	000066-25-1	38
3		Acetic acid, chloro-, butyl este...	150	C6H11ClO2	000590-02-3	10
4		Butanal, 2-methyl- (CAS) \$\$ 2-Me...	86	C5H10O	000096-17-3	9



Library Search Compound Report

Data File : C:\MSDCHEM\1\5973N\02MAR01\03184.D

Acq On : 1 Mar 2002 12:48 pm

Sample : 508 scan

Misc : March 1, 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)

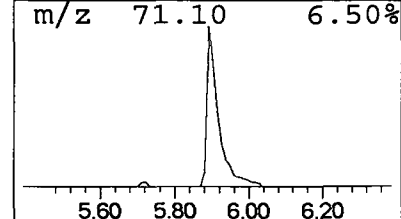
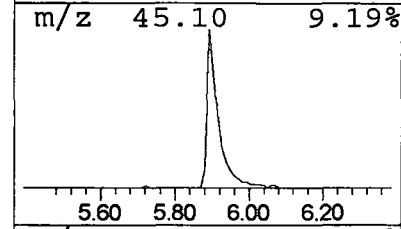
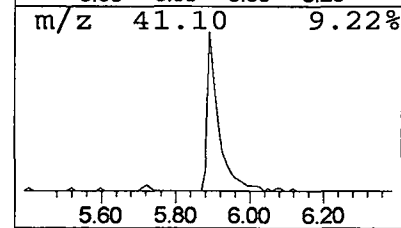
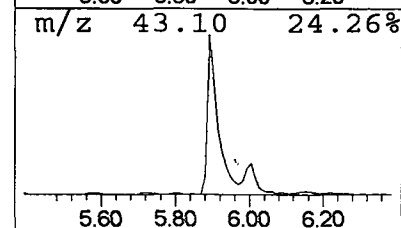
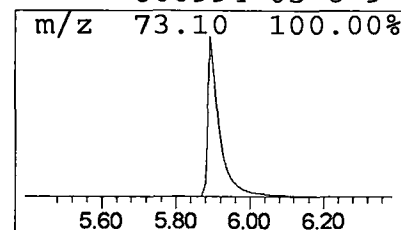
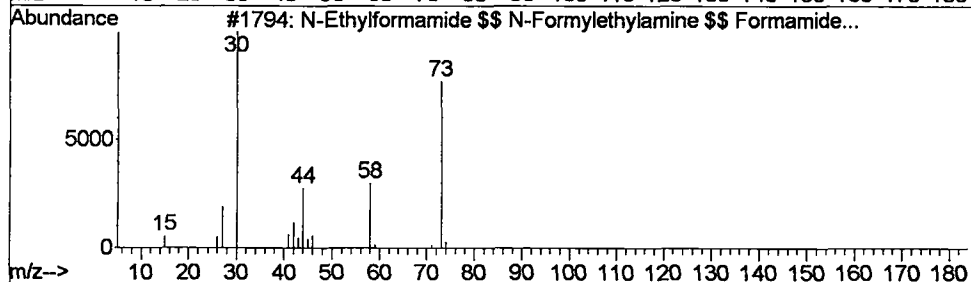
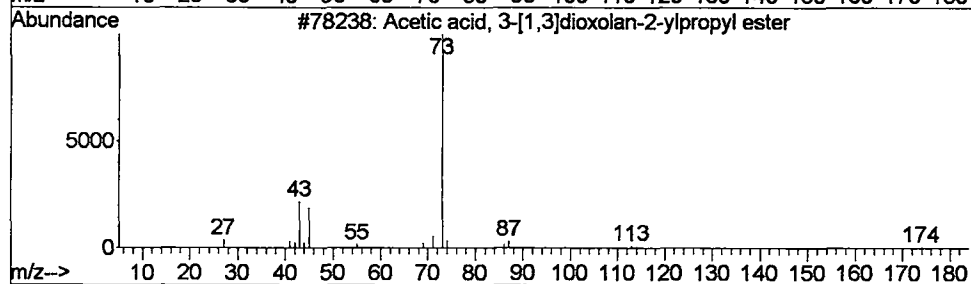
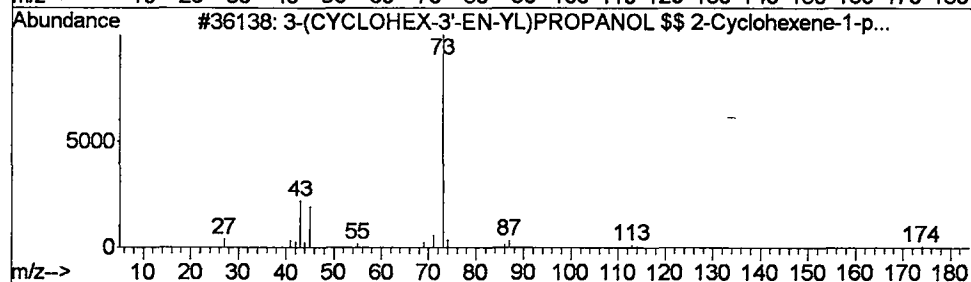
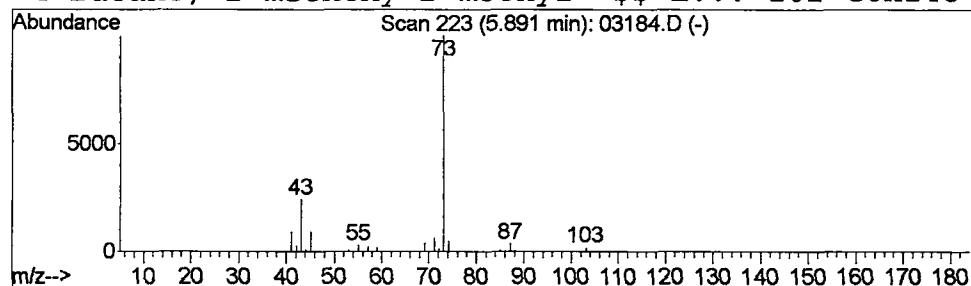
Title : 508 scan method

Library : C:\DATABASE\WILEY7N.L

Peak Number 9 3-(CYCLOHEX-3'-EN-YL)PROPAN... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.89	4.13 ug/L	1462430	1,4-Dichlorobenzene-d4	9.71

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			3-(CYCLOHEX-3'-EN-YL)PROPANOL	174	C8H14O4	015745-87-6	33
2			Acetic acid, 3-[1,3]dioxolan-2-y...	174	C8H14O4	000000-00-0	33
3			N-Ethylformamide	73	C3H7NO	000627-45-2	9
4			Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	9



Data File : C:\MSDCHEM\1\5973N\02MAR01\03184.D
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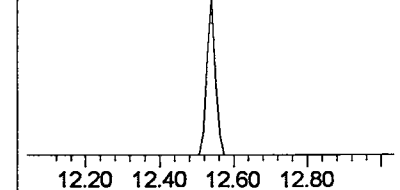
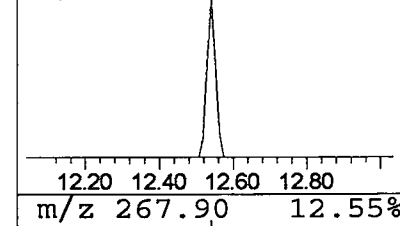
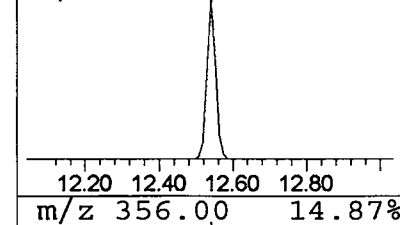
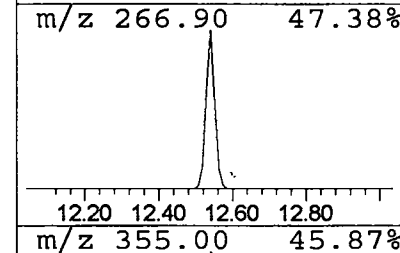
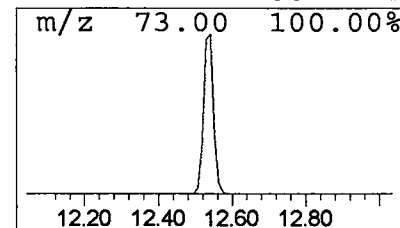
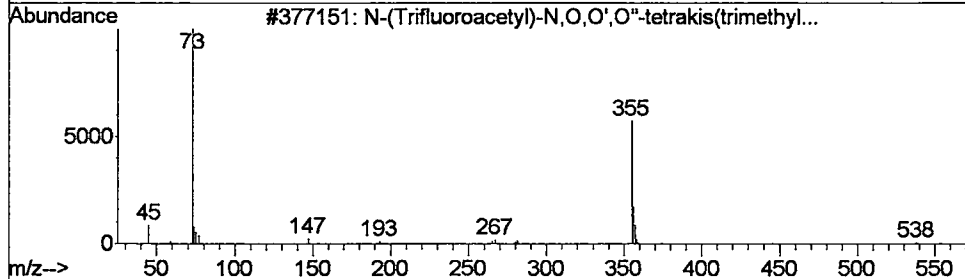
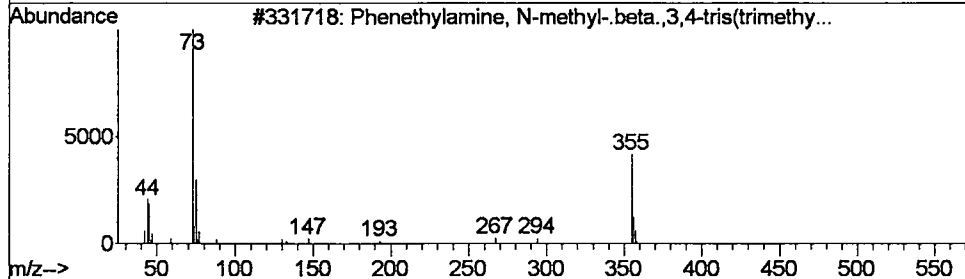
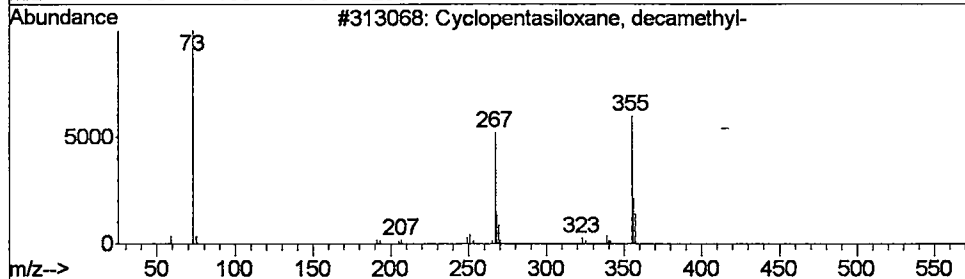
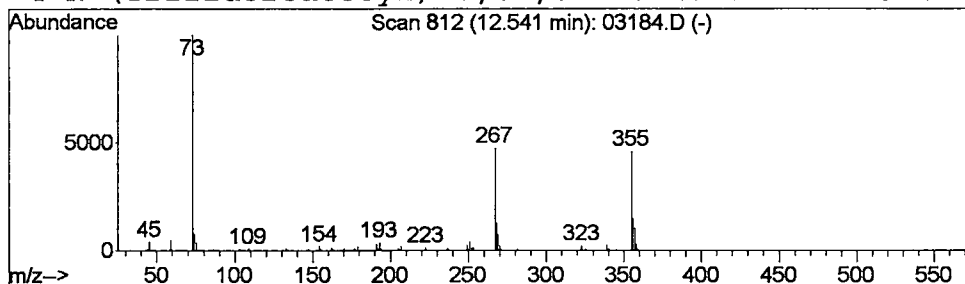
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Quant Method : C:\MSDCHEM\1...\HP022102.M (RTE Integrator)
 Title : 508 scan method
 Library : C:\DATABASE\WILEY7N.L

 Peak Number 10 Cyclopentasiloxane, decamet... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.54	0.92 ug/L	472050	Naphthalene-d8	13.20

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Cyclopentasiloxane, decamethyl-	370	C10H30O5Si5	000541-02-6	68
2		Phenethylamine, N-methyl-.beta.,...	399	C18H37NO3Si3	010538-85-9	47
3		N-(Trifluoroacetyl)-N,O,O',O''-t...	553	C22H42F3NO4Si4	000000-00-0	47
4		N-(Trifluoroacetyl)-O,O',O''-tri...	481	C19H34F3NO4Si3	000000-00-0	38



Library Search Compound Report

Data File : C:\MSDCHEM\1\5973N\02MAR01\03184.D
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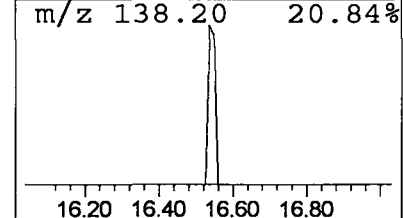
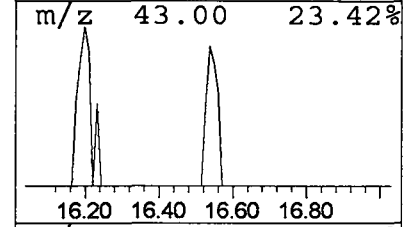
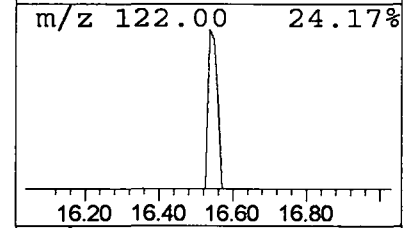
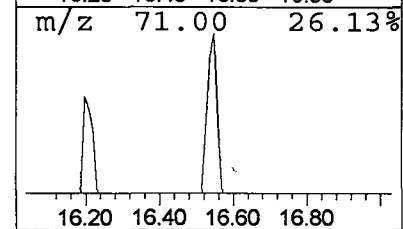
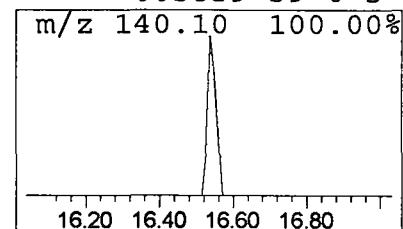
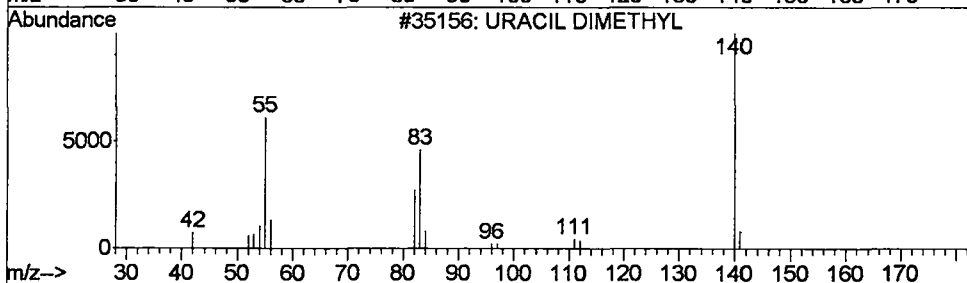
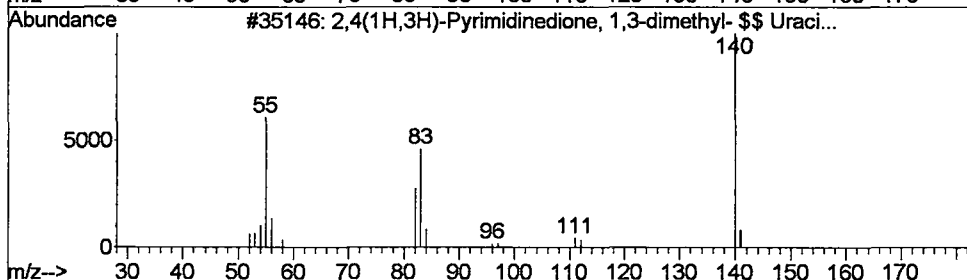
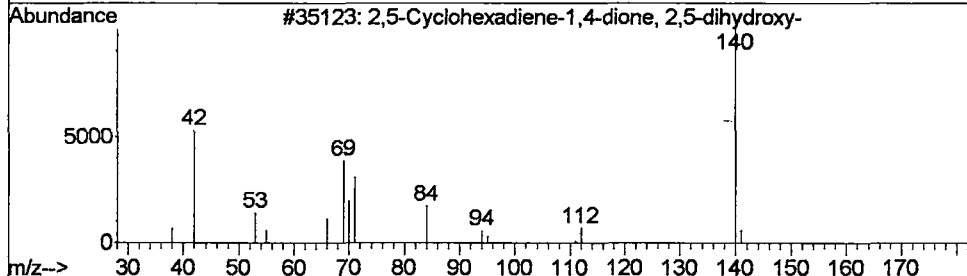
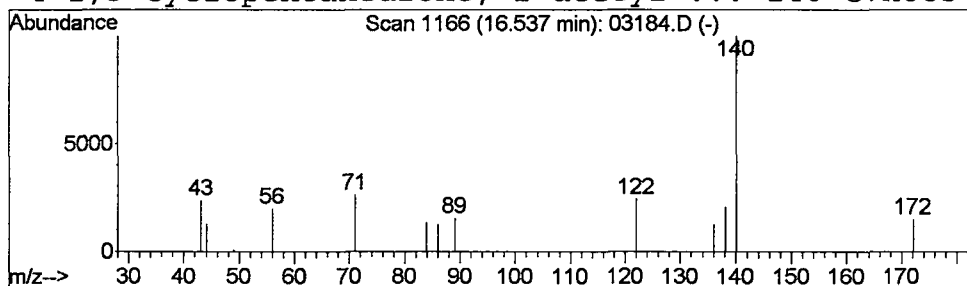
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Inst : Instrumen
Multiplr: 1.00

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

Peak Number 11 2,5-Cyclohexadiene-1,4-dione... Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.54	0.04 ug/L	25028	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			2,4(1H,3H)-Pyrimidinedione, 1,3-...	140	C6H8N2O2	000874-14-6	7
3			URACIL DIMETHYL	140	C6H8N2O2	000000-00-0	7
4			1,3-Cyclopentanedione, 2-acetyl-...	140	C7H8O3	003859-39-0	5



Library Search Compound Report

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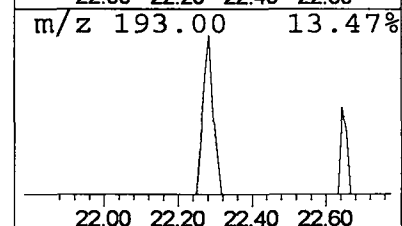
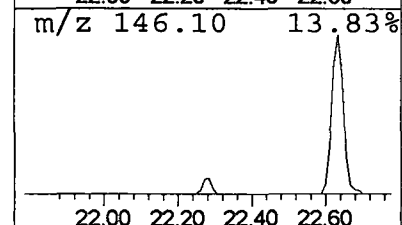
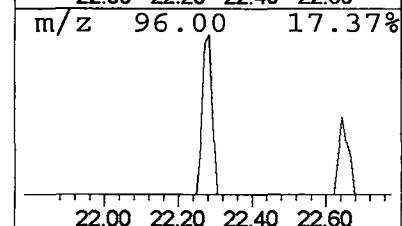
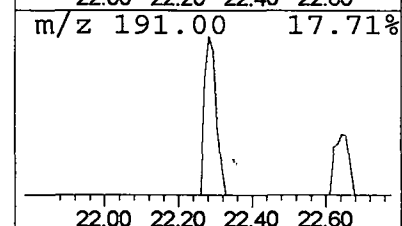
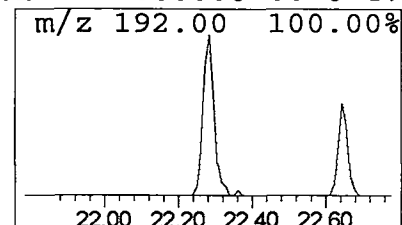
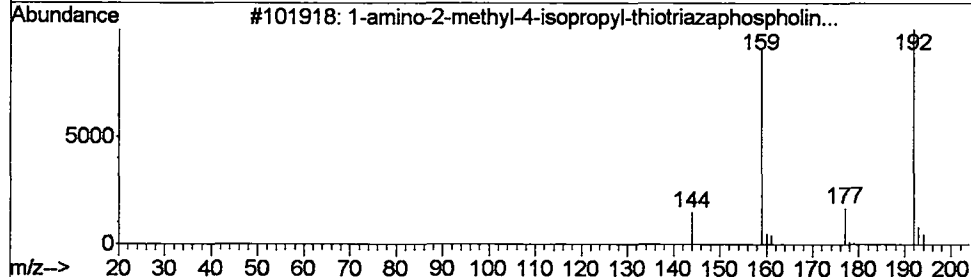
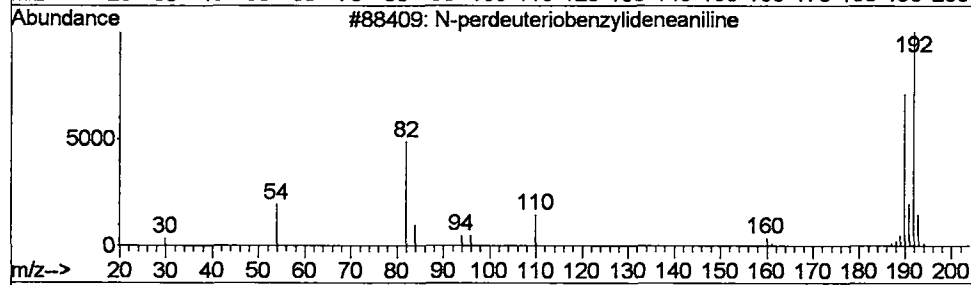
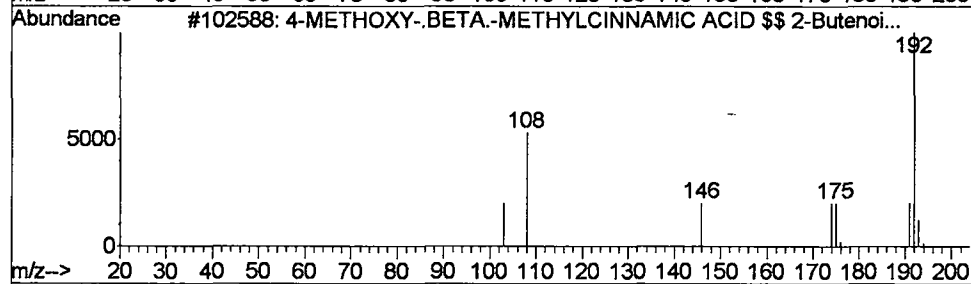
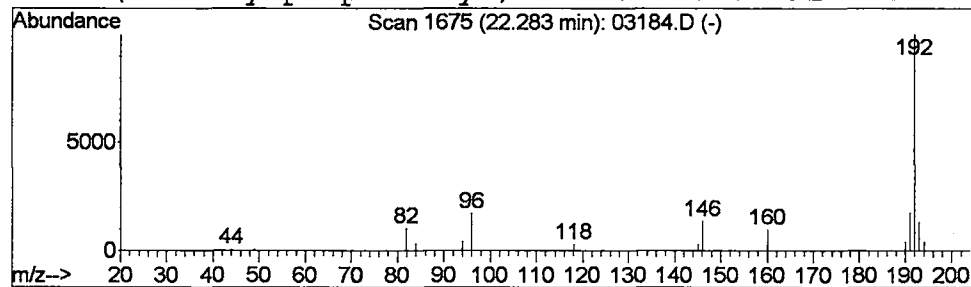
Vial: 4
Operator:
Inst : Instrumen
Multiplr: 1.00

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

Peak Number 12 4-METHOXY-.BETA.-METHYLCINN... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.28	0.12 ug/L	72108	Phenanthrene-d10	22.63

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			4-METHOXY-.BETA.-METHYLCINNAMIC ...	192	C11H12O3	019169-94-9	72
2			N-perdeuteriobenzylideneaniline	192	C13D11N	000538-51-2	53
3			1-amino-2-methyl-4-isopropyl-thi...	192	C5H13N4PS	107264-54-0	50
4			N-(2-methylprop-2-enyl)-2-nitrob...	192	C10H12N2O2	000000-00-0	47



Library Search Compound Report

Data File : C:\MSDCHEM\1\5973N\02MAR01\03184.D
 Acq On : 1 Mar 2002 12:48 pm
 Sample : 508 scan
 Misc : March 1, 2002
 MS Integration Params: LSCINT.P

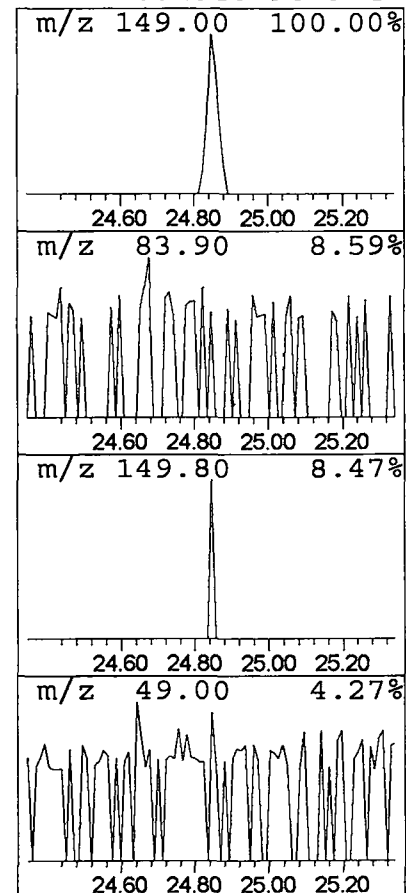
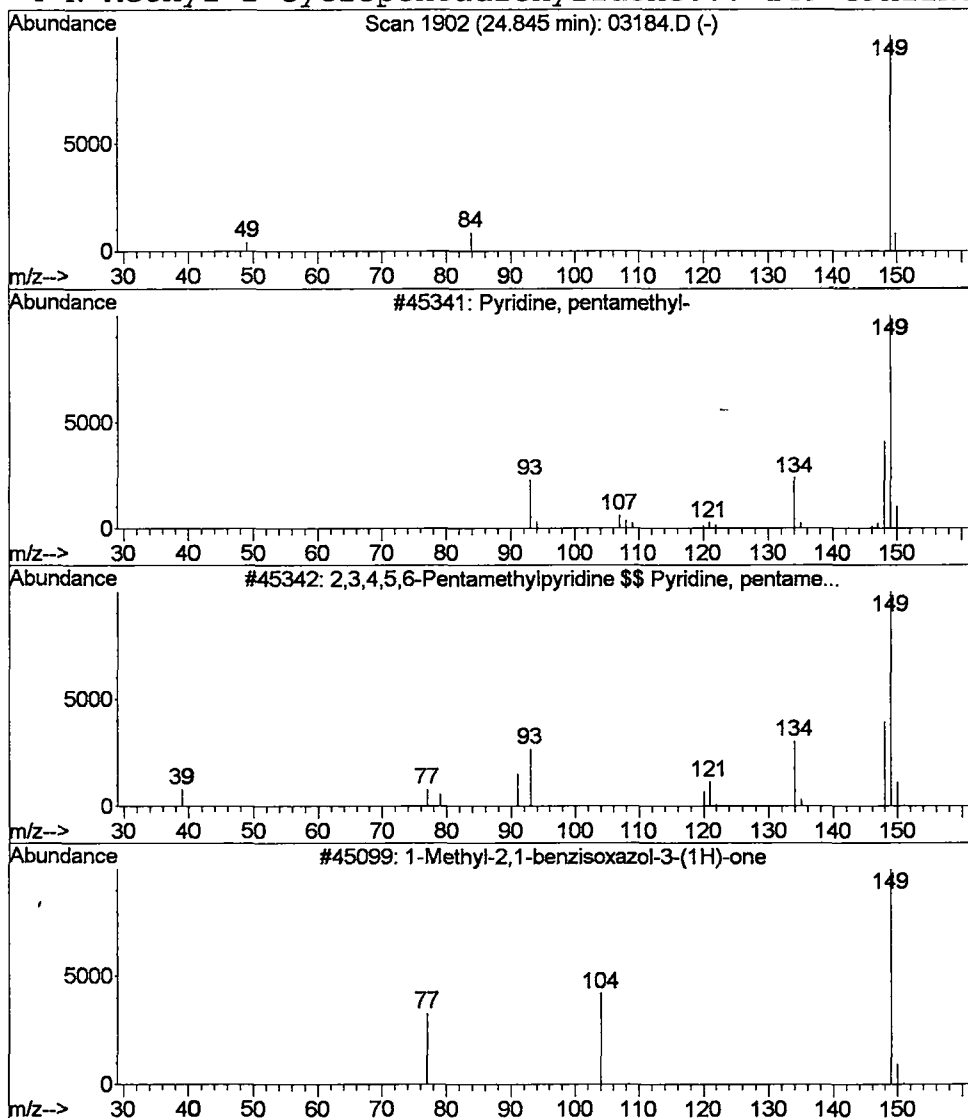
Vial: 4
 Operator:
 Inst : Instrumen
 Multiplr: 1.00

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

 Peak Number 13 Pyridine, pentamethyl- Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.85	0.03 ug/L	19955	Phenanthrene-d10	22.63

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Pyridine, pentamethyl-	149	C10H15N	003748-83-2	4
2		2,3,4,5,6-Pentamethylpyridine \$\$...	149	C10H15N	003748-83-2	4
3		1-Methyl-2,1-benzisoxazol-3-(1H)...	149	C8H7NO2	000000-00-0	4
4		N-Methyl-2-cyclopentadienylidene...	149	C9H11NO	081540-98-9	4



tentatively identified compound (LSC) summary

Operator ID: Date Acquired: 1 Mar 2002 12:48 pm
 Data File: C:\MSDCHEM\1\5973N\02MAR01\03184.D
 Name: 508 scan
 Date: March 1, 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
Acetonitrile, dic...	3.49	0.1	ug/L	24096	1	9.71	7077220	20.0
Butanoic acid, me...	3.62	0.1	ug/L	51419	1	9.71	7077220	20.0
Propane, 1,1-dime...	3.93	0.1	ug/L	24714	1	9.71	7077220	20.0
1,4-DI-TRIDEUTERO...	4.27	0.1	ug/L	21157	1	9.71	7077220	20.0
2-Buten-1-ol, 3-m...	4.65	0.1	ug/L	38457	1	9.71	7077220	20.0
2,3,4,5-Tetrahydr...	4.86	0.2	ug/L	81939	1	9.71	7077220	20.0
3utenol, methyl-	5.02	0.1	ug/L	21857	1	9.71	7077220	20.0
Hexanal (CAS) \$\$...	5.11	0.1	ug/L	20556	1	9.71	7077220	20.0
3-(CYCLOHEX-3'-EN...	5.89	4.1	ug/L	1462430	1	9.71	7077220	20.0
Cyclopentasiloxan...	12.54	0.9	ug/L	472050	2	13.20	10269000	20.0
2,5-Cyclohexadien...	16.54	0.0	ug/L	25028	3	18.33	11215300	20.0
4-METHOXY-.BETA.-...	22.28	0.1	ug/L	72108	4	22.63	12186300	20.0
Pyridine, pentame...	24.85	0.0	ug/L	19955	4	22.63	12186300	20.0