

User: Vinci, David L
 Westchester Dept. of Labs & Research
 Date: 03/05/2002 Time: 16:53:43

Sample: AE01510
 Analysis: \$GCMS GCMS Scan For Unknowns
 Units: ug
 Started: 3/5/02 16:52 Ended: 3/5/02 16:52 Analyst: DLV
 Page: 1

	Component Name	Viol	Result	Qualifier	MDL
1	Other unknown compounds		.		
2	Surr_2,4-DDD		92		5.0

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· **Comments for Sample AE01510 - Analysis \$GCMS**

Westchester Dept. of Labs & Research

User: Vinci, David L

Date: 03-05-2002 Time: 16:53:47

The GCMS scan, from 35 to 550 amu, does not reveal the presence of unusual compounds. The recoveries for 22 of the 43 compounds in the LCS Standard were above their acceptance ranges.

Quantitation Report

(Not Reviewed)

Data File : C:\MSDCHEM\1\5973N\02FEB21\01510.D

Vial: 5

Acq On : 21 Feb 2002 2:17 pm

Operator:

Sample : 508 scan

Inst : Instrumen

Misc : February 21, 2002

Multiplr: 1.00

MS Integration Params: SPLIT.P

Quant Time: Feb 25 10:49:09 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	1643471	20.00	ug/L	0.00
10) Naphthalene-d8	13.20	136	7210128	20.00	ug/L	0.00
17) Acenaphthene-d10	18.34	164	3639932	20.00	ug/L	0.00
29) Phenanthrene-d10	22.64	188	6301621	20.00	ug/L	0.00
38) Chrysene-d12	30.49	240	4580140	20.00	ug/L	-0.03
44) Perylene-d12	34.41	264	3059438	20.00	ug/L	-0.04

System Monitoring Compounds

37) 2,4-DDD	27.72	235	387117	4.61	ug/L	0.00
Spiked Amount	5.000		Recovery	=	92.20%	

Target Compounds

						Qvalue
20) Dimethylphthalate	18.34	163	584140	2.65	ug/L	# 58
21) 2,6-Dinitrotoluene	18.34	165	456422	8.95	ug/L	# 27

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

01510.D HP022102.M Mon Feb 25 10:49:10 2002

Page 1

LSC Area Percent Report

Data File : C:\MSDCHEM\1\5973N\02FEB21\01510.D

Acq On : 21 Feb 2002 2:17 pm

Sample : 508 scan

Misc : February 21, 2002

MS Integration Params: LSCINT.P

Vial: 5

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	3.498	9	11	19	rVB	25816	42630	0.28%	0.053%
2	3.622	19	22	29	rVB	44340	70261	0.46%	0.087%
3	3.927	47	49	55	rVB	22984	37728	0.24%	0.047%
4	4.378	85	89	95	rVB	21246	32697	0.21%	0.041%
5	5.948	225	228	237	rBV2	64643	240299	1.56%	0.298%
6	6.060	237	238	255	rVB	69730	227971	1.48%	0.283%
7	9.718	557	562	567	rBV	5026342	9133219	59.16%	11.323%
8	9.876	573	576	593	rVB	15818	77760	0.50%	0.096%
9	13.195	865	870	875	rBV	7049848	13611437	88.16%	16.876%
10	13.376	883	886	893	rVB2	14432	32628	0.21%	0.040%
11	18.343	1321	1326	1331	rBV	7332585	14701118	95.22%	18.227%
12	19.980	1467	1471	1481	rVB3	17562	51884	0.34%	0.064%
13	22.283	1671	1675	1683	rVB	48329	98114	0.64%	0.122%
14	22.644	1701	1707	1711	rBV	7004077	15439152	100.00%	19.142%
15	22.768	1715	1718	1741	rVB	103193	529173	3.43%	0.656%
16	24.856	1899	1903	1909	rBV	15906	33952	0.22%	0.042%
17	27.724	2153	2157	2163	rVV	1036914	1888779	12.23%	2.342%
18	30.490	2397	2402	2441	rBV2	5589942	13744739	89.03%	17.041%
19	33.402	2655	2660	2665	rBV2	45364	100968	0.65%	0.125%
20	34.418	2743	2750	2779	rBV	3640342	10563245	68.42%	13.096%

Sum of corrected areas: 80657754

Quantitation Report

(Not Reviewed)

Data File : C:\MSDCHEM\1\5973N\02FEB21\01510.D

Vial: 5

Acq On : 21 Feb 2002 2:17 pm

Operator:

Sample : 508 scan

Inst : Instrumen

Misc : February 21, 2002

Multiplr: 1.00

MS Integration Params: SPLIT.P

Quant Time: Feb 25 10:49 2002

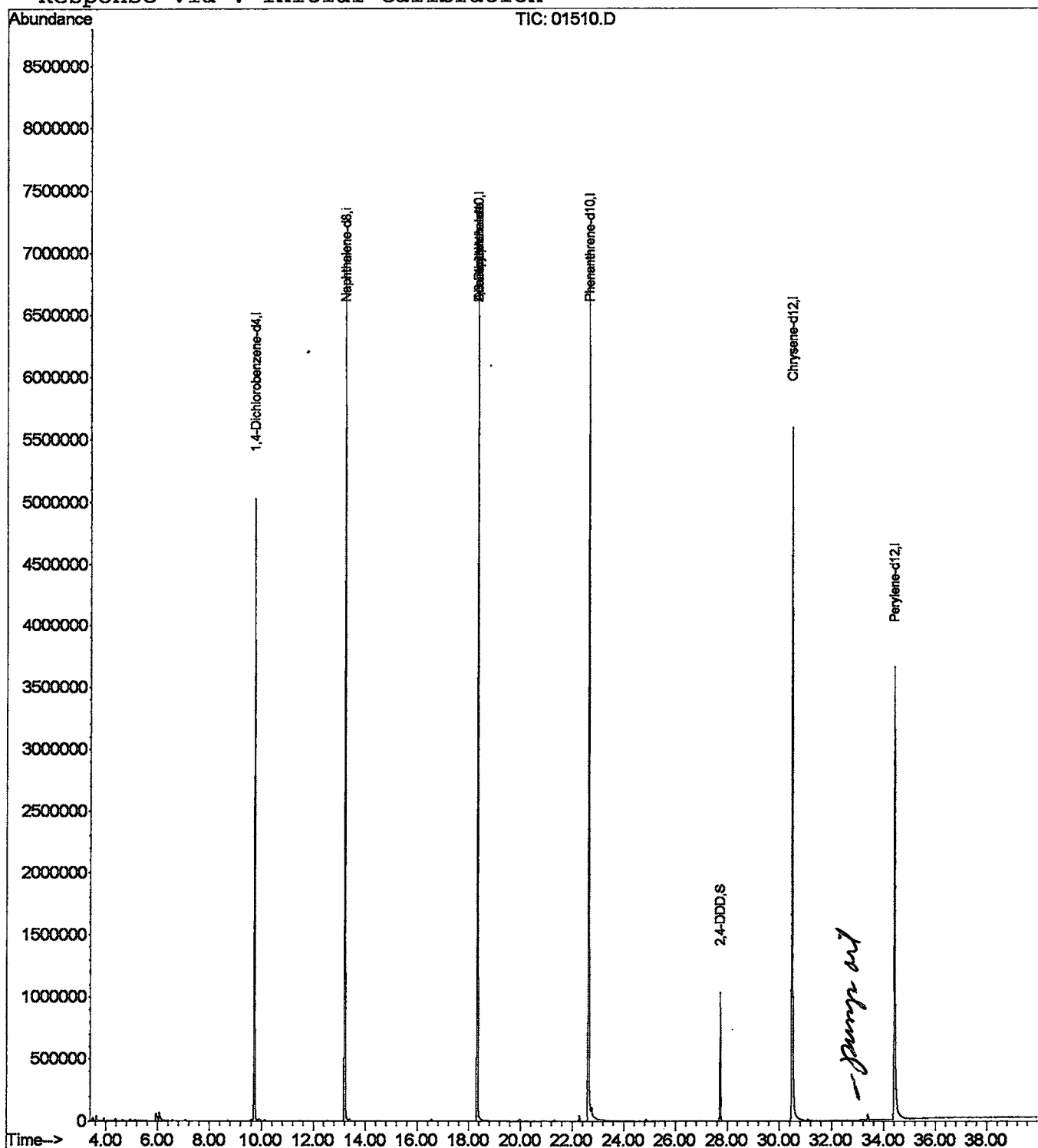
Quant Results File: HP022102.R

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

Title : 508 scan method

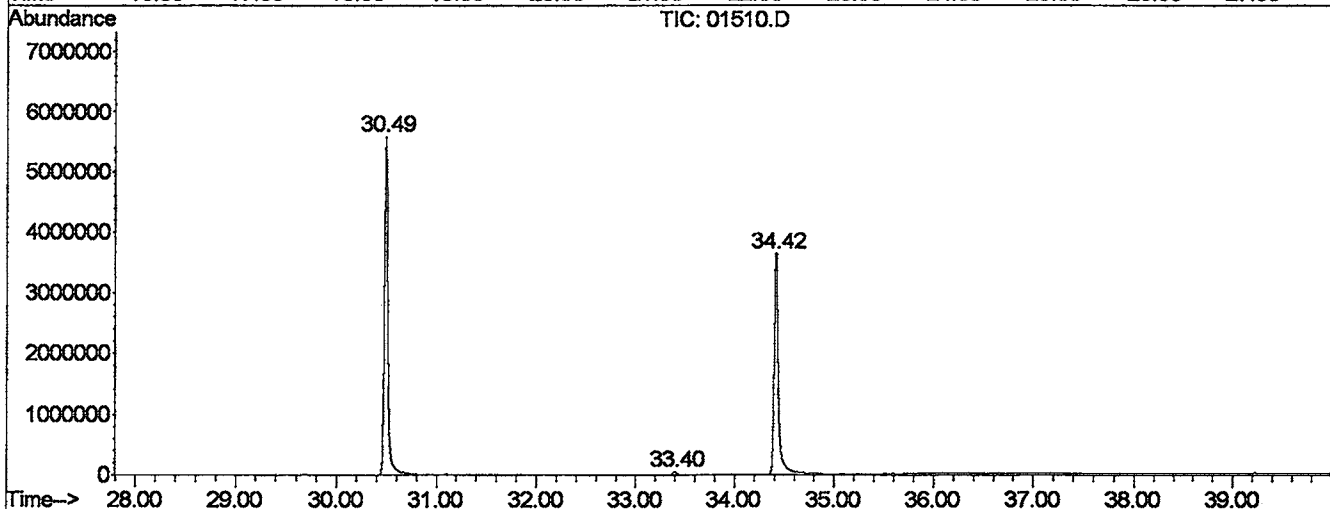
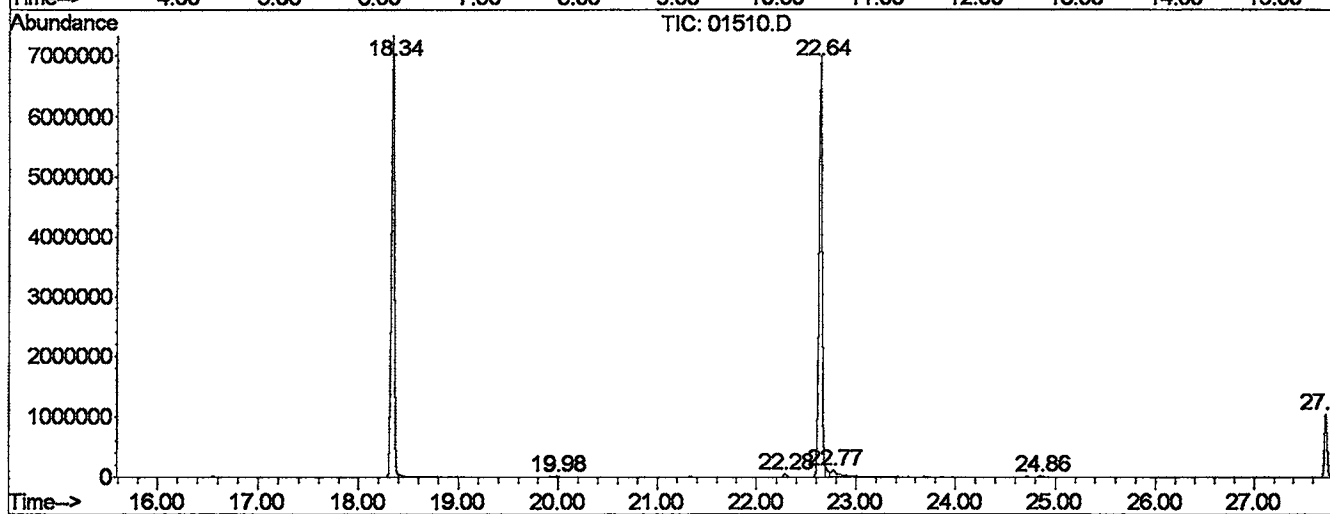
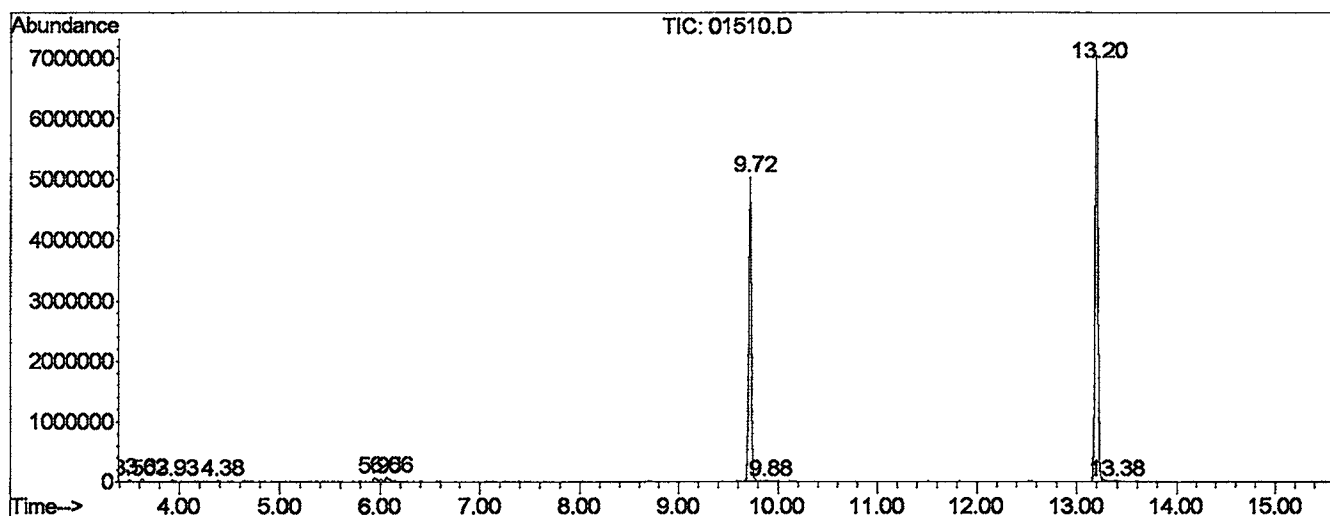
Last Update : Mon Feb 25 10:47:59 2002

Response via : Initial Calibration



LSC Report - Integrated Chromatogram

File : C:\MSDCHEM\1\5973N\02FEB21\01510.D
 Operator :
 Acquired : 21 Feb 2002 2:17 pm using AcqMethod HP020702
 Instrument : Instrumen
 Sample Name: 508 scan
 Misc Info : February 21, 2002
 Vial Number: 5
 Quant File :HP022102.RES (RTE Integrator)



Library Search Compound Report

Data File : C:\MSDCHEM\1\5973N\02FEB21\01510.D
 Acq On : 21 Feb 2002 2:17 pm
 Sample : 508 scan
 Misc : February 21, 2002
 MS Integration Params: LSCINT.P

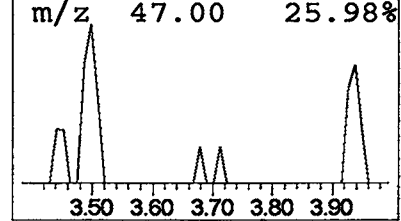
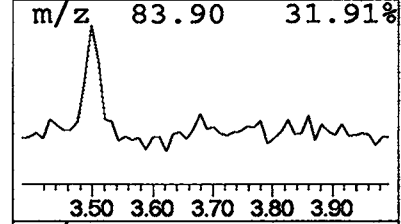
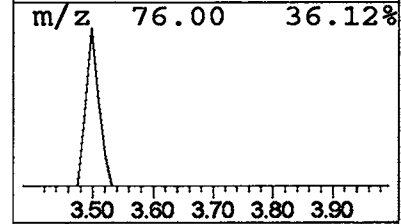
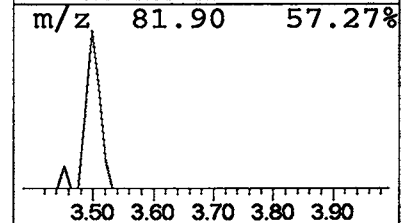
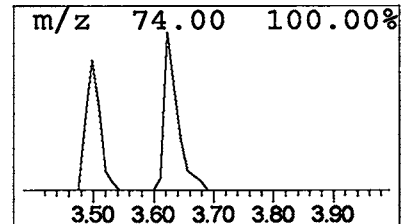
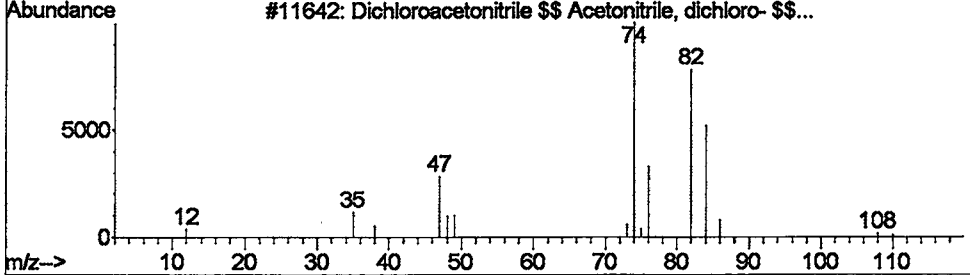
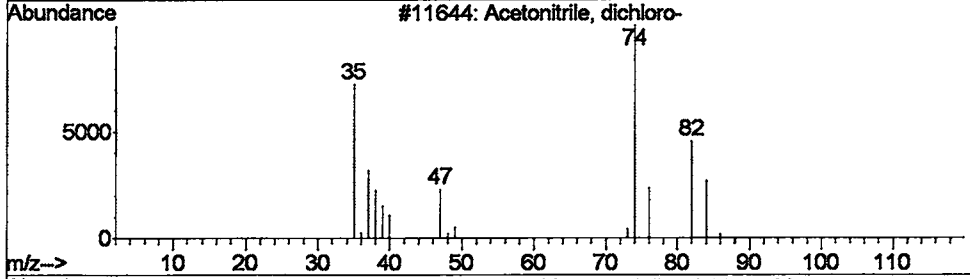
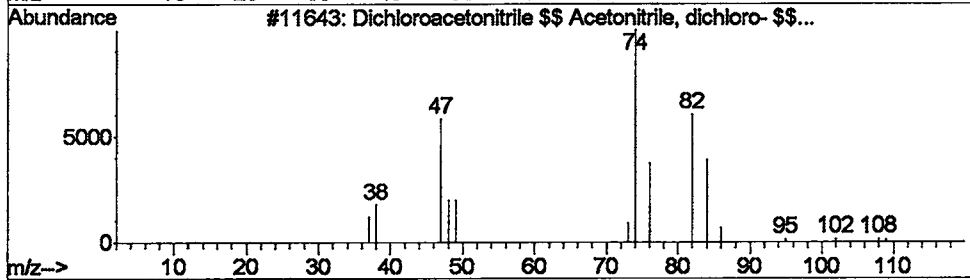
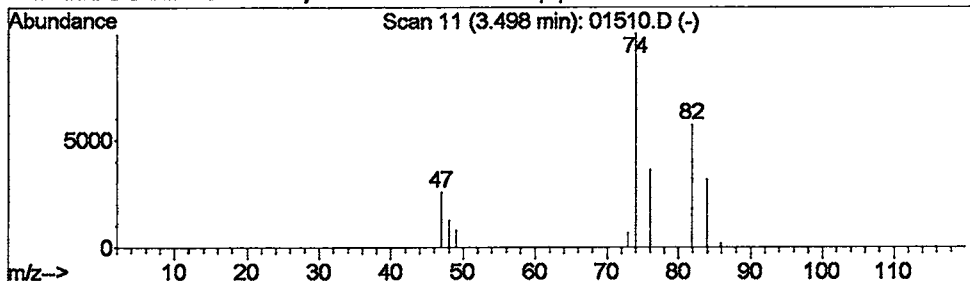
Vial: 5
 Operator:
 Inst : Instrumen
 Multiplr: 1.00

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

 Peak Number 1 Dichloroacetonitrile \$\$ Ace... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.50	0.09 ug/L	42630	1,4-Dichlorobenzene-d4	9.72

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Dichloroacetonitrile \$\$ Acetonit...	109	C2HCl2N	003018-12-0	83
2			Acetonitrile, dichloro-	109	C2HCl2N	003018-12-0	83
3			Dichloroacetonitrile \$\$ Acetonit...	109	C2HCl2N	003018-12-0	74
4			Acetonitrile, dichloro- \$\$ Dichl...	109	C2HCl2N	003018-12-0	74



Library Search Compound Report

Data File : C:\MSDCHEM\1\5973N\02FEB21\01510.D
 Acq On : 21 Feb 2002 2:17 pm
 Sample : 508 scan
 Misc : February 21, 2002
 MS Integration Params: LSCINT.P

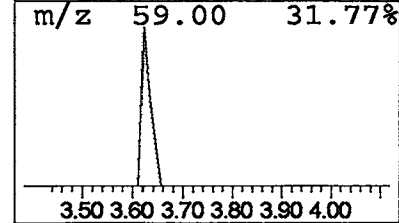
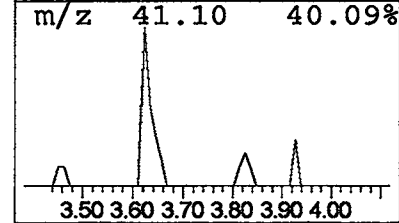
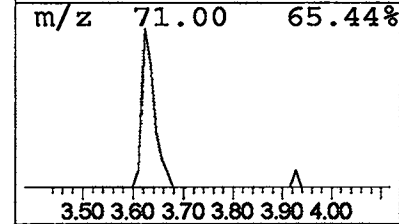
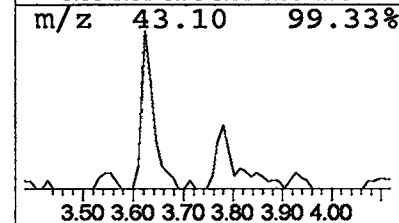
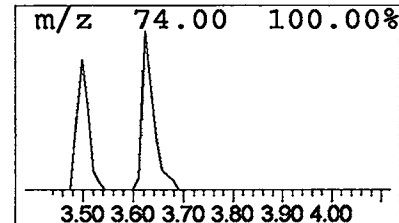
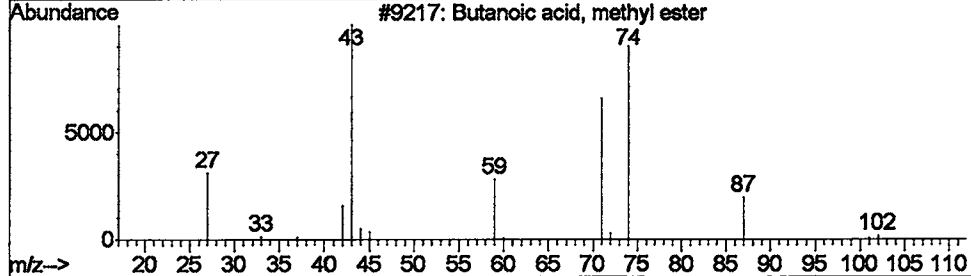
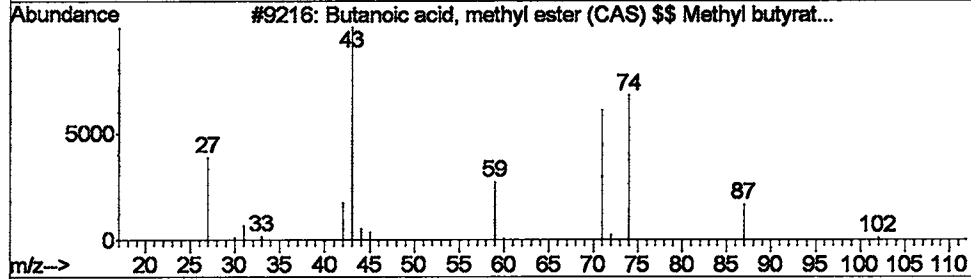
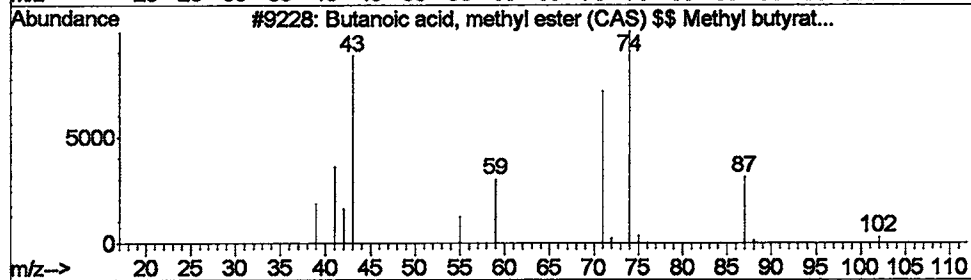
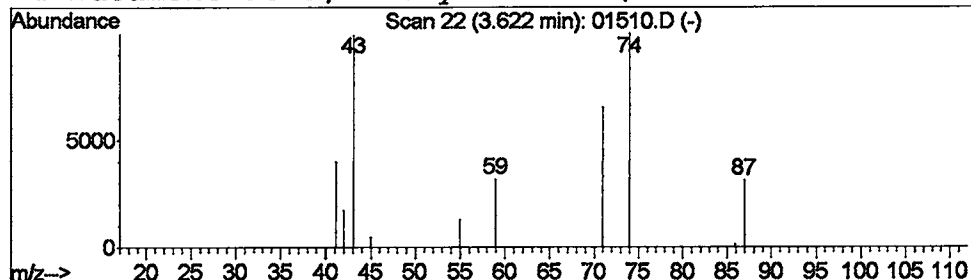
Vial: 5
 Operator:
 Inst : Instrumen
 Multiplr: 1.00

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 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.62	0.15 ug/L	70261	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 (CAS...	102	C5H10O2	000623-42-7	83
3			Butanoic acid, methyl ester	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\02FEB21\01510.D
Acq On : 21 Feb 2002 2:17 pm
Sample : 508 scan
Misc : February 21, 2002
MS Integration Params: LSCINT.P

Vial: 5
Operator:
Inst : Instrumen
Multiplr: 1.00

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

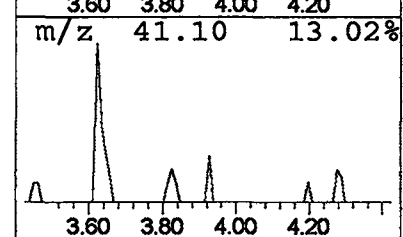
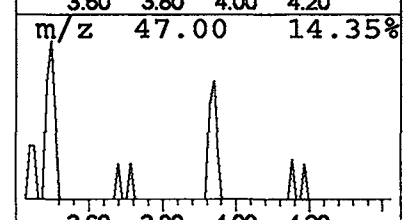
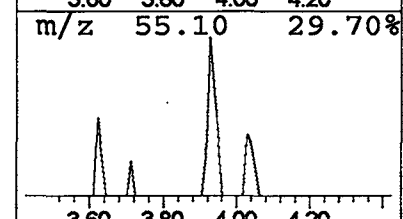
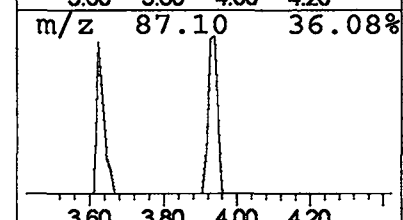
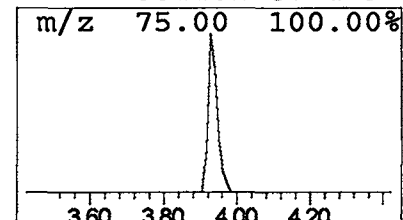
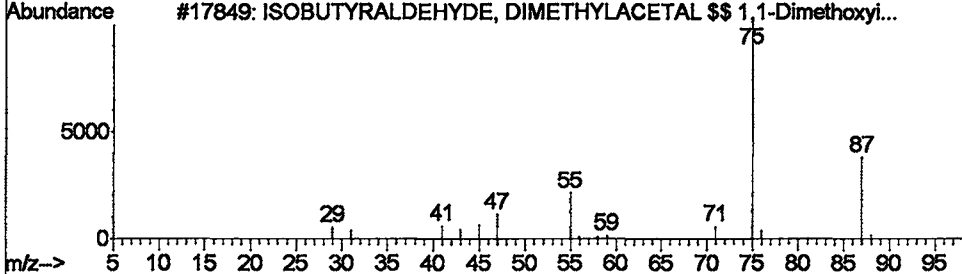
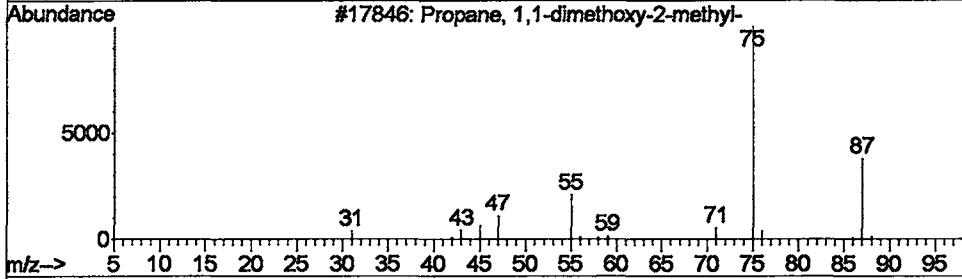
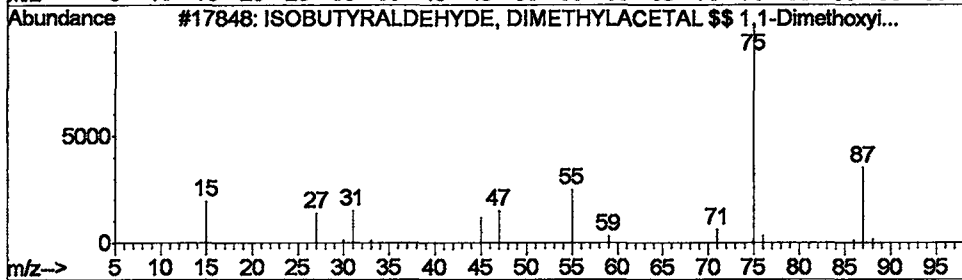
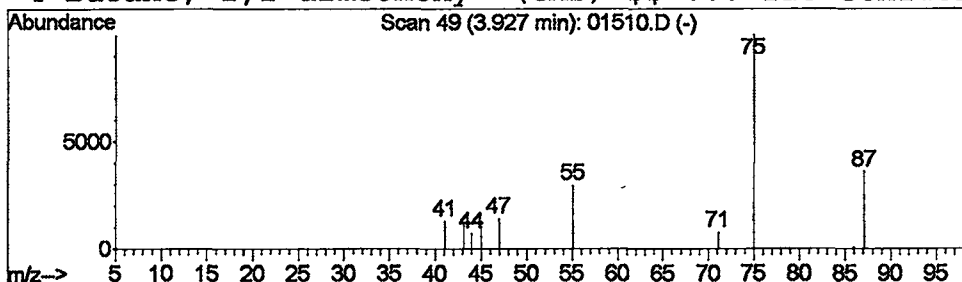
Title : 508 scan method

Library : C:\DATABASE\WILEY7N.L

Peak Number 3 ISOBUTYRALDEHYDE, DIMETHYLA... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.93	0.08 ug/L	37728	1,4-Dichlorobenzene-d4	9.72

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			ISOBUTYRALDEHYDE, DIMETHYLACETAL...	118	C6H14O2	000000-00-0	83
2			Propane, 1,1-dimethoxy-2-methyl-	118	C6H14O2	041632-89-7	83
3			ISOBUTYRALDEHYDE, DIMETHYLACETAL...	118	C6H14O2	000000-00-0	83
4			Butane, 1,1-dimethoxy- (CAS) \$\$...	118	C6H14O2	004461-87-4	9



Library Search Compound Report

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 Sample : 508 scan
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 MS Integration Params: LSCINT.P

Vial: 5
 Operator:
 Inst : Instrumen
 Multiplr: 1.00

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

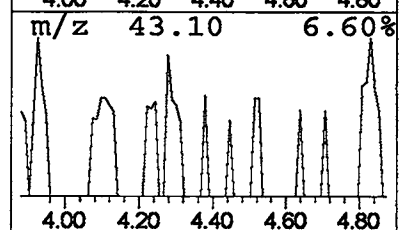
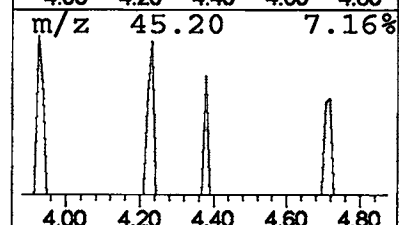
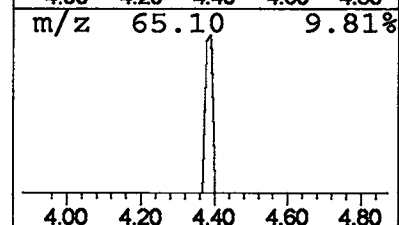
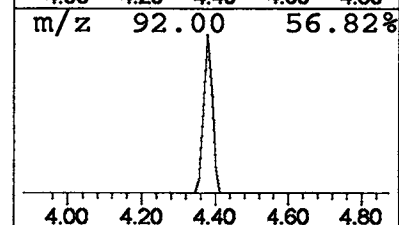
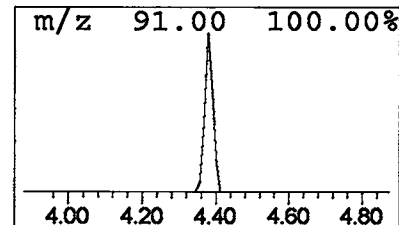
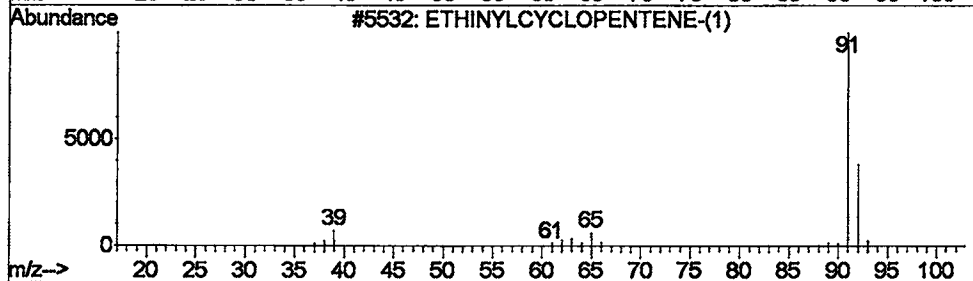
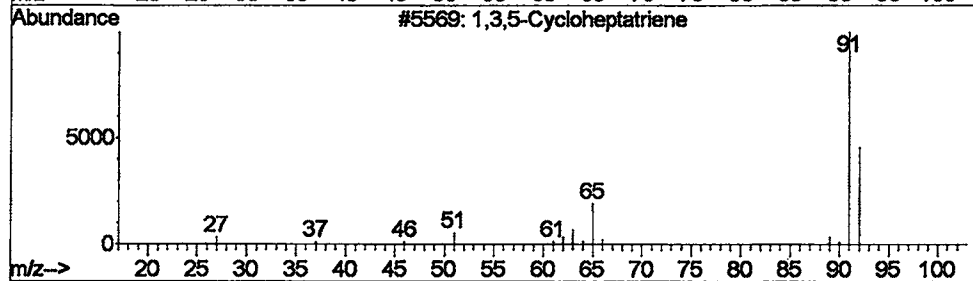
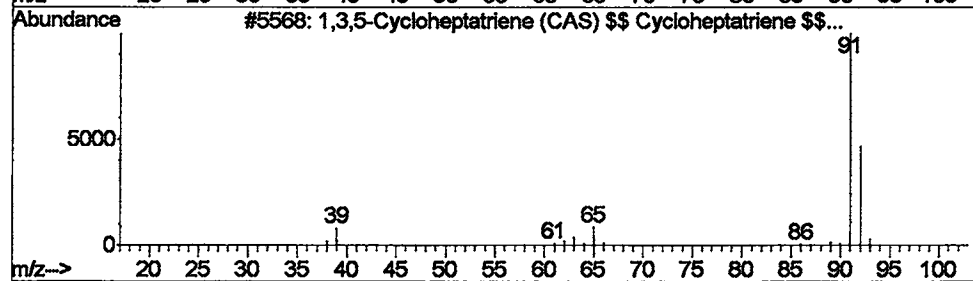
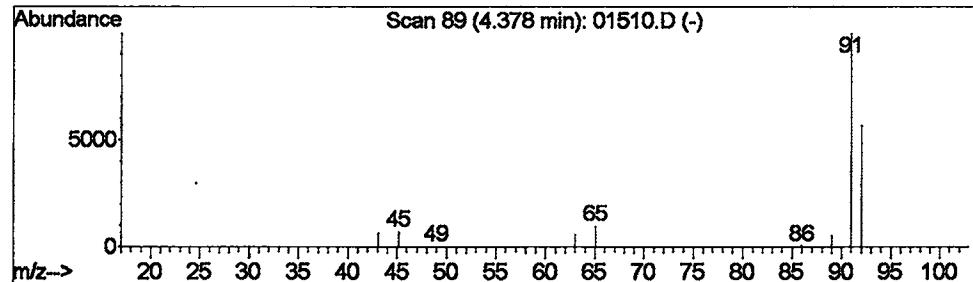
Title : 508 scan method

Library : C:\DATABASE\WILEY7N.L

 Peak Number 4 1,3,5-Cycloheptatriene (CAS... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.38	0.07 ug/L	32697	1,4-Dichlorobenzene-d4	9.72

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1,3,5-Cycloheptatriene (CAS) \$\$...	92	C7H8	000544-25-2	64
2			1,3,5-Cycloheptatriene	92	C7H8	000544-25-2	40
3			ETHINYLCYCLOPENTENE-(1)	92	C7H8	000000-00-0	9
4			METHYLFULVENE	92	C7H8	000000-00-0	9



Library Search Compound Report

Data File : C:\MSDCHEM\1\5973N\02FEB21\01510.D
Acq On : 21 Feb 2002 2:17 pm
Sample : 508 scan
Misc : February 21, 2002
MS Integration Params: LSCINT.P

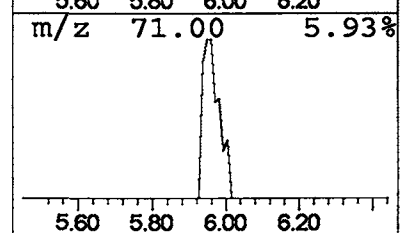
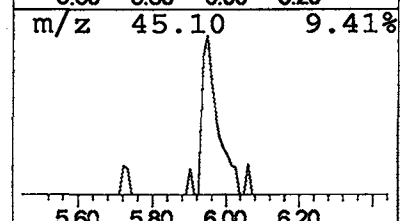
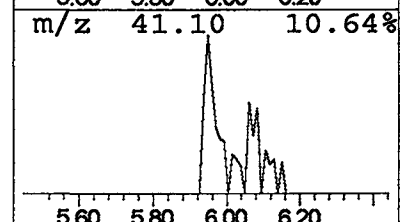
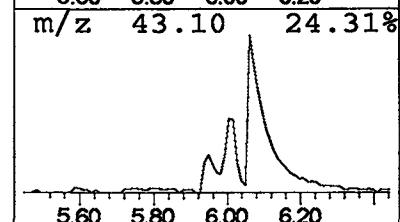
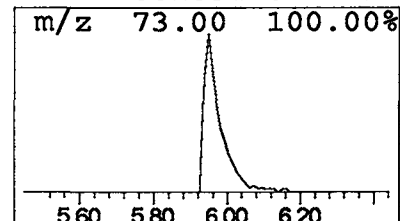
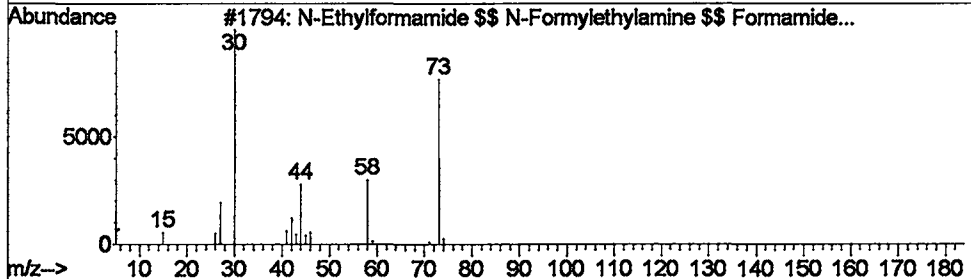
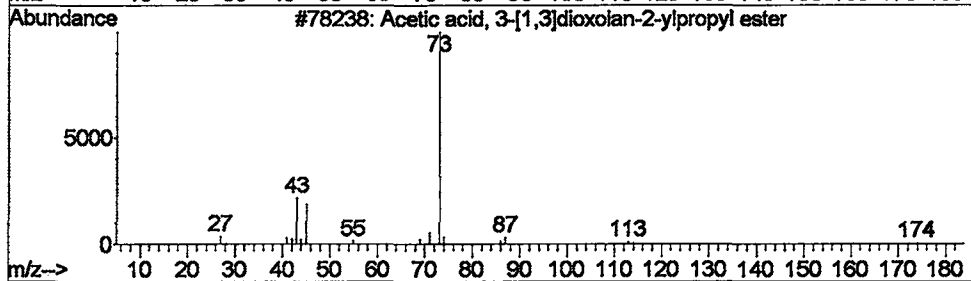
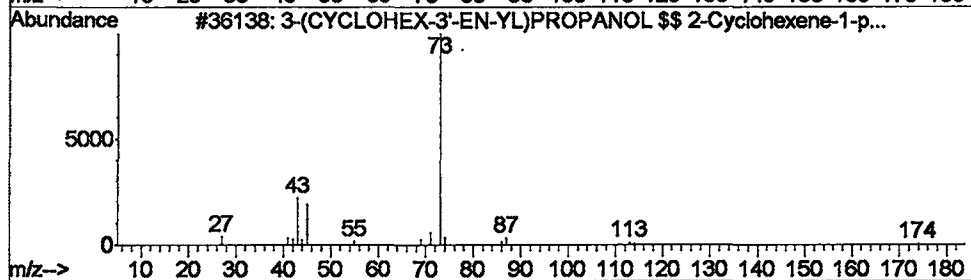
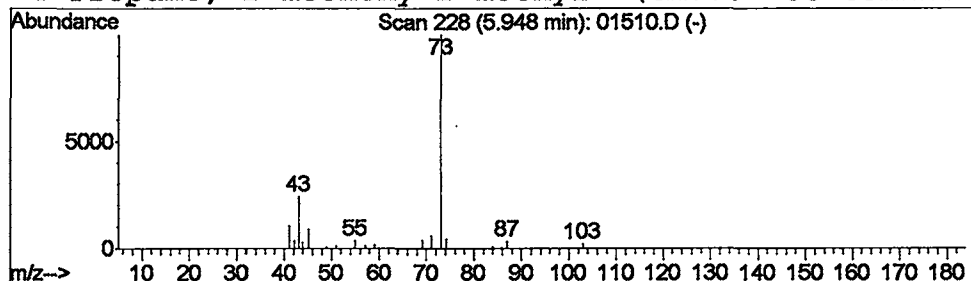
Vial: 5
Operator:
Inst : Instrumen
Multiplr: 1.00

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

Peak Number 5 3-(CYCLOHEX-3'-EN-YL)PROPAN... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.95	0.53 ug/L	240299	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			Propane, 2-methoxy-2-methyl- (CA...	88	C5H12O	001634-04-4	9



Library Search Compound Report

Data File : C:\MSDCHEM\1\5973N\02FEB21\01510.D
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 Misc : February 21, 2002
 MS Integration Params: LSCINT.P

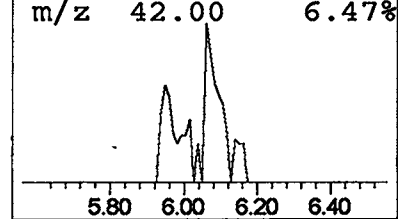
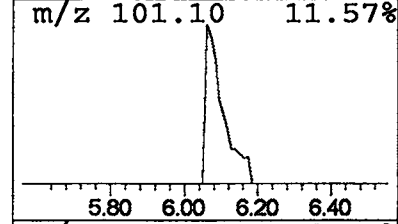
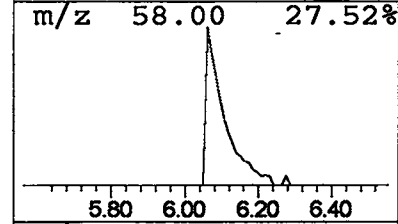
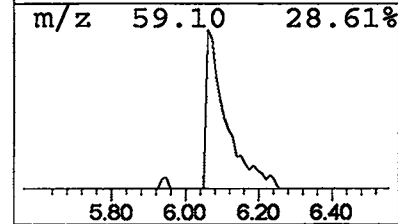
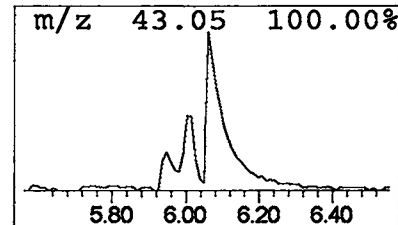
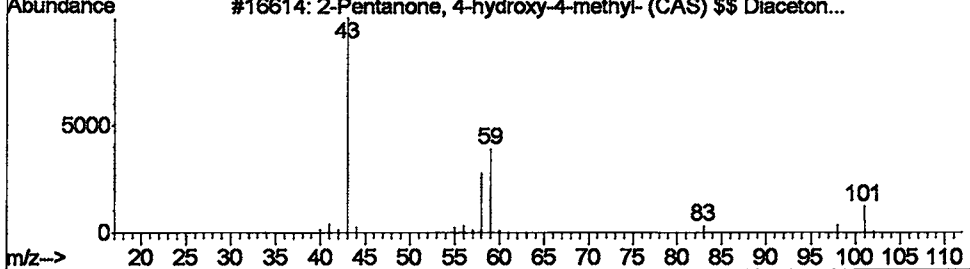
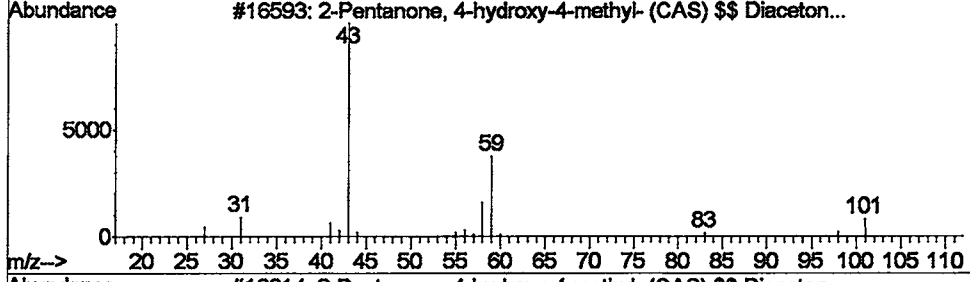
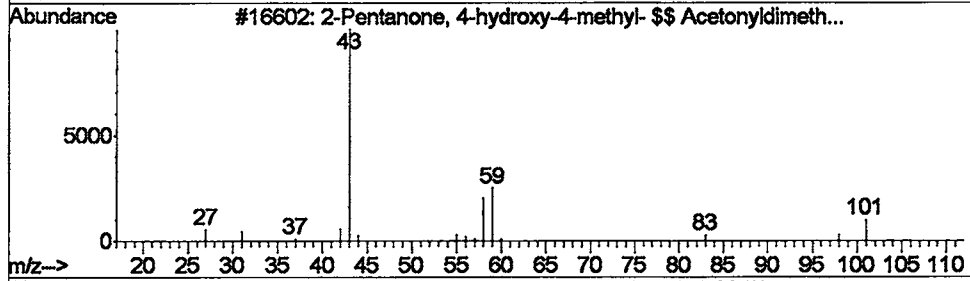
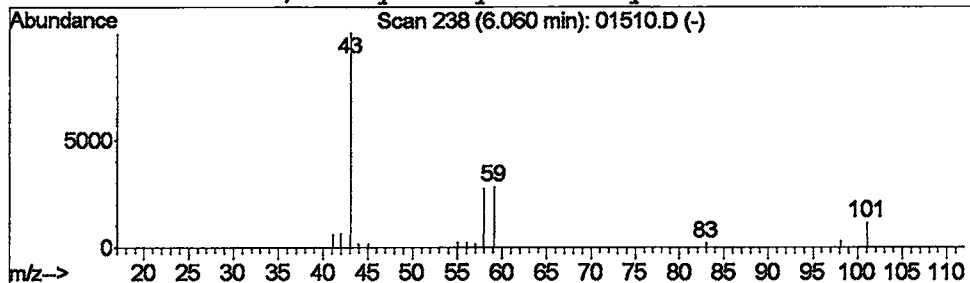
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 Inst : Instrumen
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Quant Method : C:\MSDCHEM\1...\HP022102.M (RTE Integrator)
 Title : 508 scan method
 Library : C:\DATABASE\WILEY7N.L

 Peak Number 6 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.06	0.50 ug/L	227971	1,4-Dichlorobenzene-d4	9.72

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentanone, 4-hydroxy-4-methyl-...	116	C6H12O2	000123-42-2	74
2			2-Pentanone, 4-hydroxy-4-methyl-...	116	C6H12O2	000123-42-2	72
3			2-Pentanone, 4-hydroxy-4-methyl-...	116	C6H12O2	000123-42-2	64
4			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	64



Library Search Compound Report

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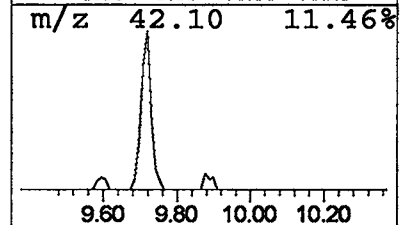
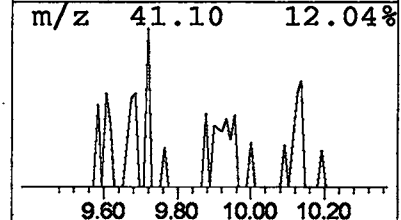
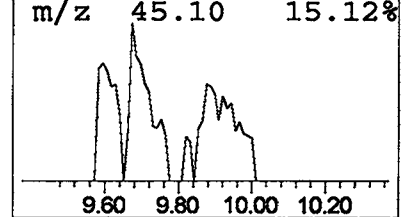
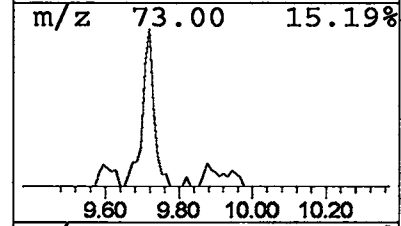
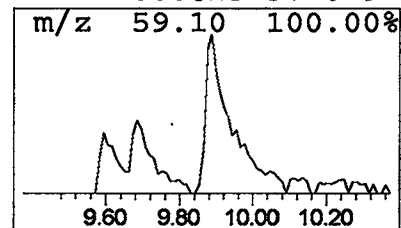
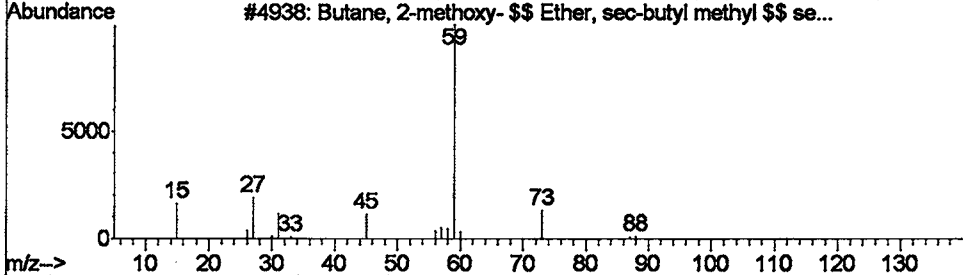
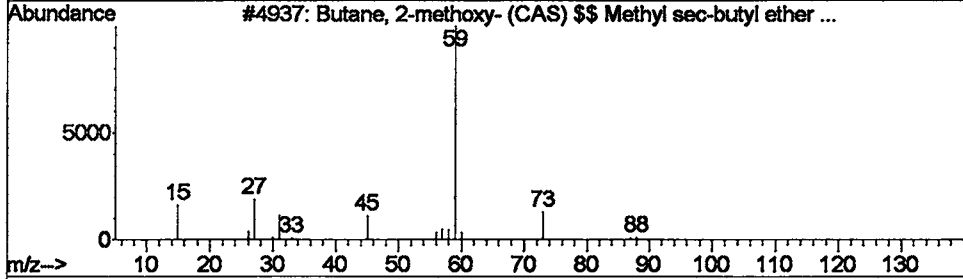
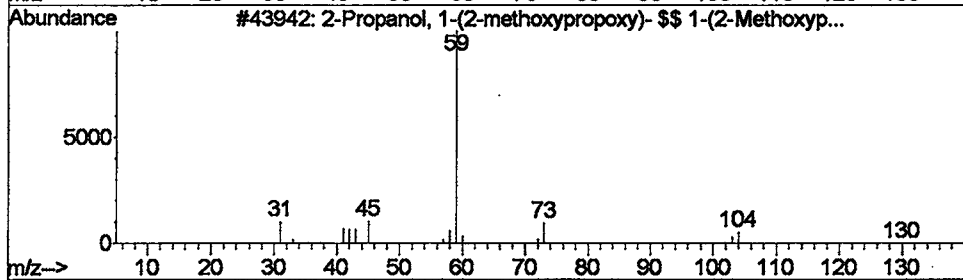
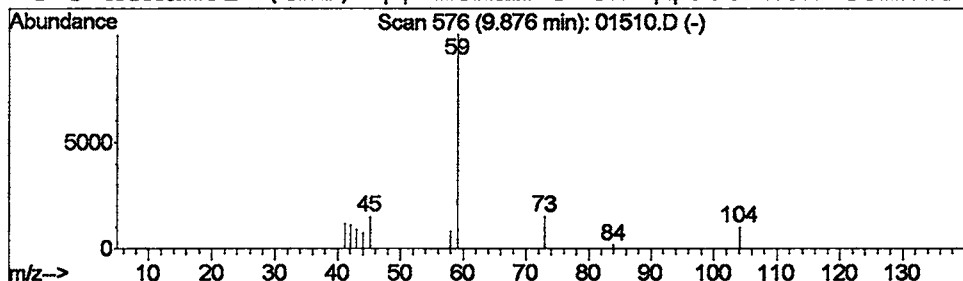
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Multiplr: 1.00

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

Peak Number 7 2-Propanol, 1-(2-methoxypro... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.88	0.17 ug/L	77760	1,4-Dichlorobenzene-d4	9.72

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Propanol, 1-(2-methoxypropoxy)...	148	C7H16O3	013429-07-7	74
2			Butane, 2-methoxy- (CAS) \$\$ Meth...	88	C5H12O	006795-87-5	9
3			Butane, 2-methoxy- \$\$ Ether, sec...	88	C5H12O	006795-87-5	9
4			3-Hexanol (CAS) \$\$ Hexan-3-ol \$\$...	102	C6H14O	000623-37-0	9



Library Search Compound Report

Data File : C:\MSDCHEM\1\5973N\02FEB21\01510.D
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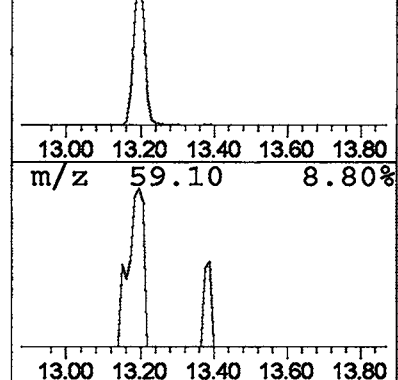
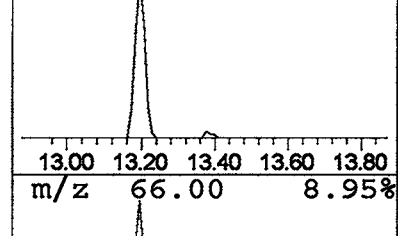
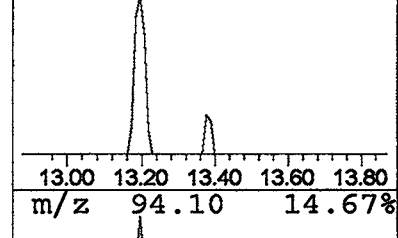
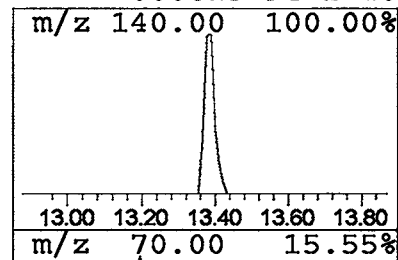
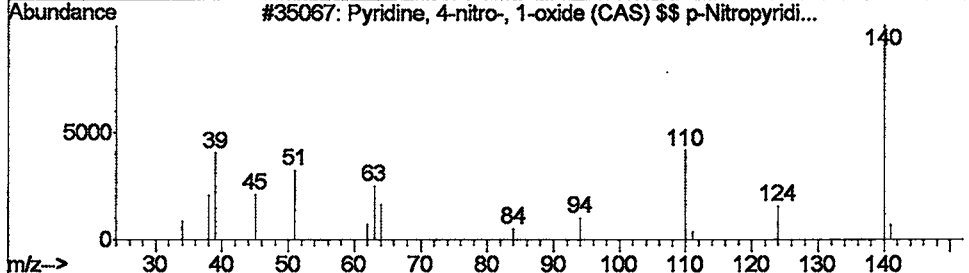
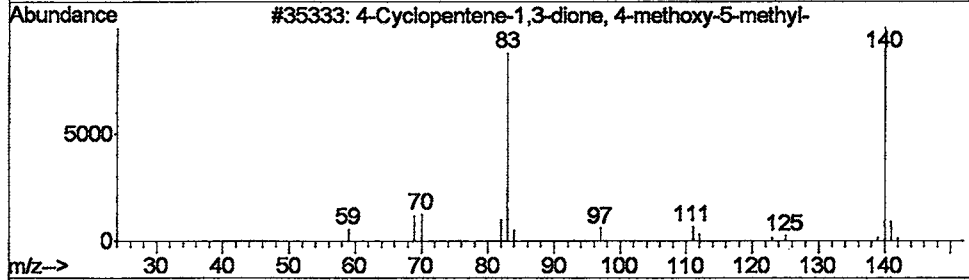
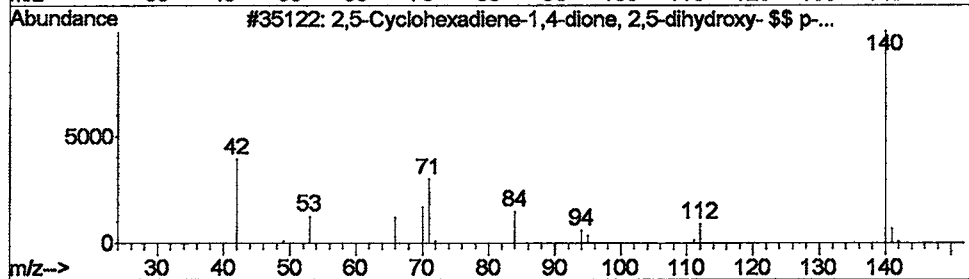
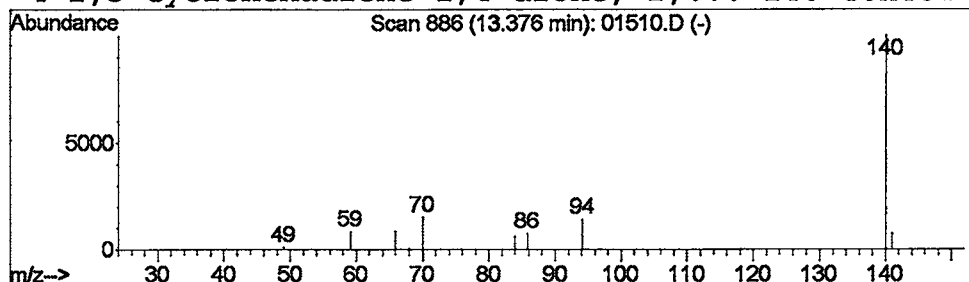
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Quant Method : C:\MSDCHEM\1...\HP022102.M (RTE Integrator)
 Title : 508 scan method
 Library : C:\DATABASE\WILEY7N.L

 Peak Number 8 2,5-Cyclohexadiene-1,4-dione... Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.38	0.05 ug/L	32628	Naphthalene-d8	13.20

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		2,5-Cyclohexadiene-1,4-dione, 2,...	140	C6H4O4	000615-94-1	72
2		4-Cyclopentene-1,3-dione, 4-meth...	140	C7H8O3	007180-62-3	64
3		Pyridine, 4-nitro-, 1-oxide (CAS...	140	C5H4N2O3	001124-33-0	50
4		2,5-Cyclohexadiene-1,4-dione, 2,...	140	C6H4O4	000615-94-1	40



Library Search Compound Report

Data File : C:\MSDCHEM\1\5973N\02FEB21\01510.D
 Acq On : 21 Feb 2002 2:17 pm
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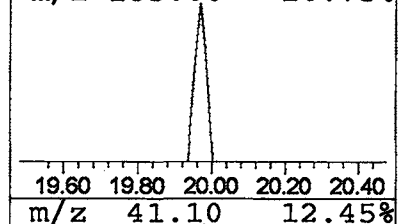
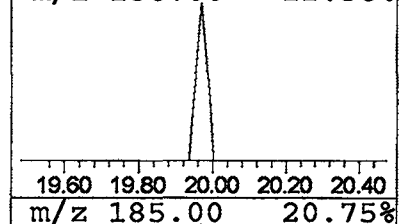
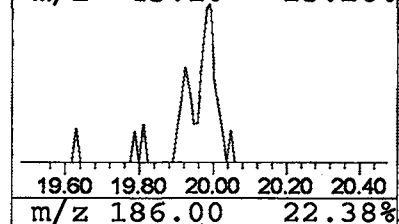
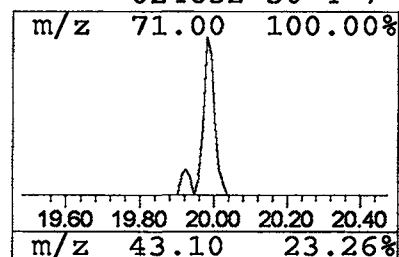
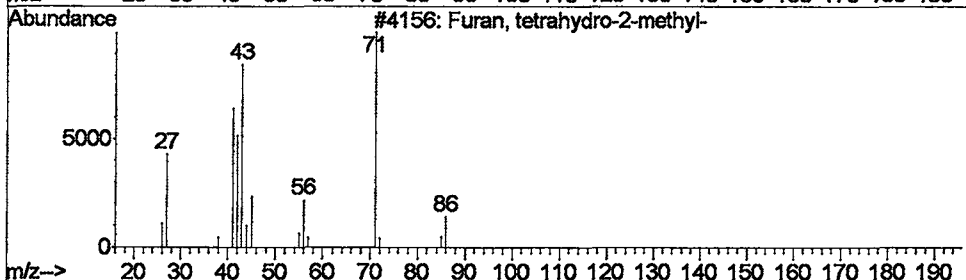
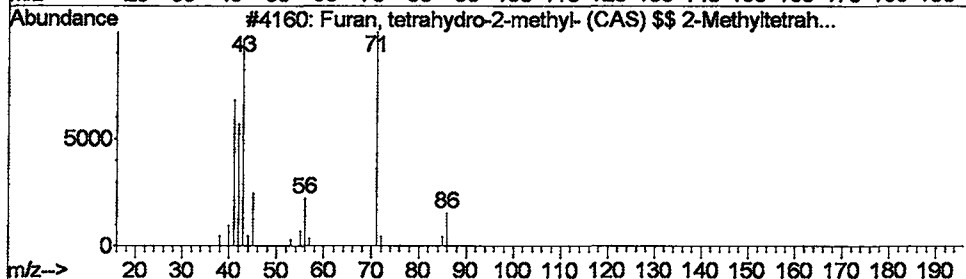
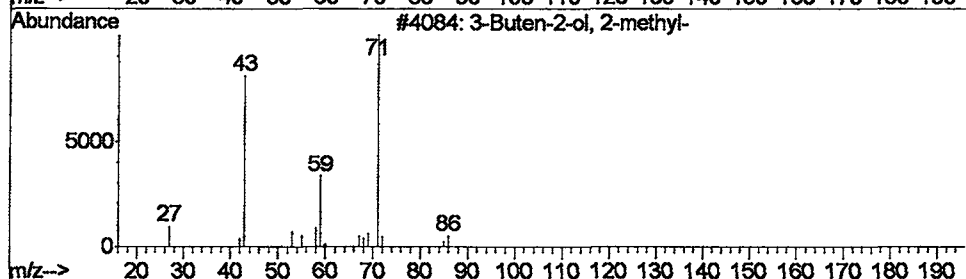
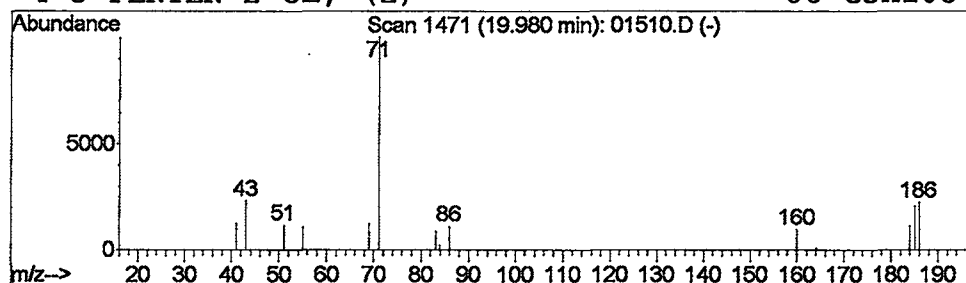
Title : 508 scan method

Library : C:\DATABASE\WILEY7N.L

 Peak Number 9 3-Buten-2-ol, 2-methyl- Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.98	0.07 ug/L	51884	Acenaphthene-d10	18.34

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		3-Buten-2-ol, 2-methyl-	86	C5H10O	000115-18-4	9
2		Furan, tetrahydro-2-methyl- (CAS...	86	C5H10O	000096-47-9	7
3		Furan, tetrahydro-2-methyl-	86	C5H10O	000096-47-9	7
4		3-PENTEN-2-OL, (Z)-	86	C5H10O	024652-50-4	7



Library Search Compound Report

Data File : C:\MSDCHEM\1\5973N\02FEB21\01510.D
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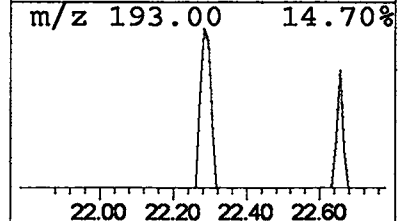
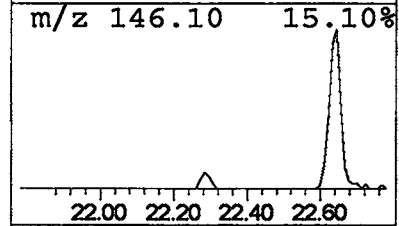
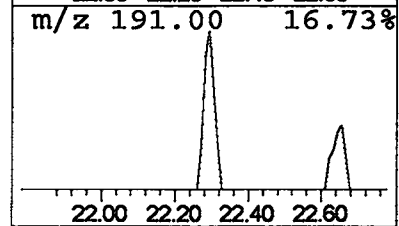
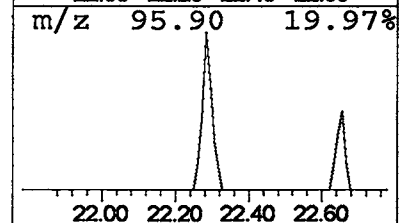
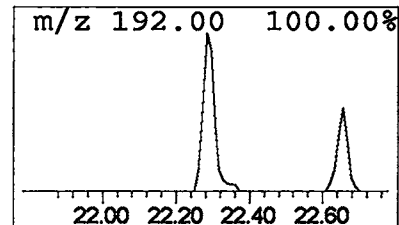
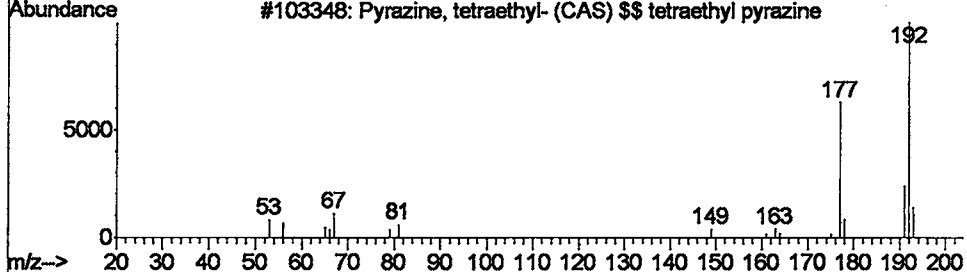
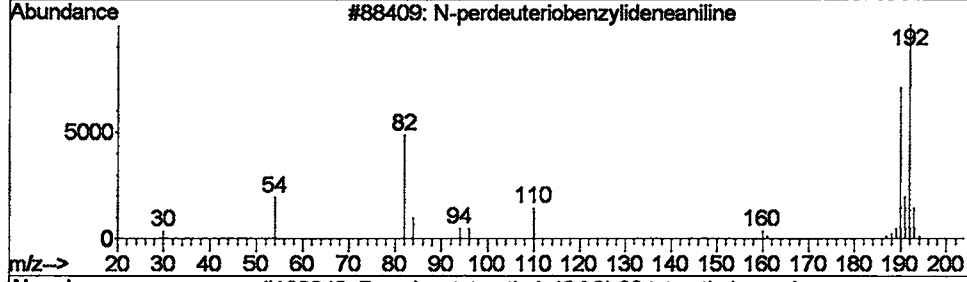
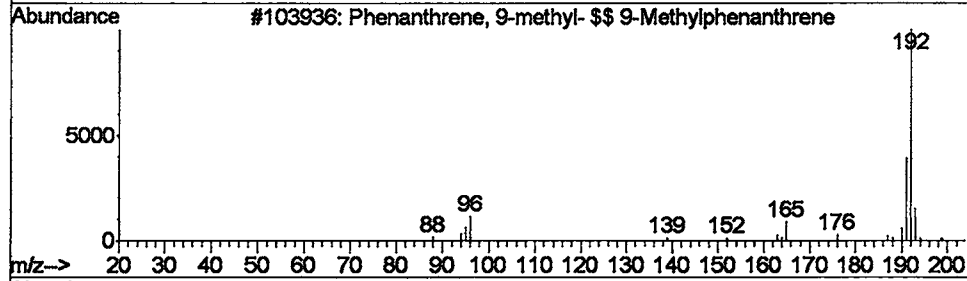
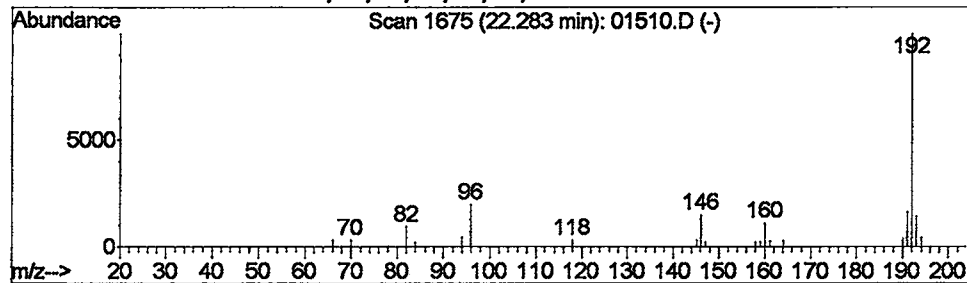
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Quant Method : C:\MSDCHEM\1...\HP022102.M (RTE Integrator)
 Title : 508 scan method
 Library : C:\DATABASE\WILEY7N.L

 Peak Number 10 Phenanthrene, 9-methyl- \$\$... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.28	0.13 ug/L	98114	Phenanthrene-d10	22.64

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Phenanthrene, 9-methyl- \$\$ 9-Met...	192	C15H12	000883-20-5	53
2			N-perdeuteriobenzylideneaniline	192	C13D11N	000538-51-2	52
3			Pyrazine, tetraethyl- (CAS) \$\$ t...	192	C12H20N2	038325-19-8	47
4			1H-2-Indenone, 2,4,5,6,7,7a-hexah...	192	C13H20O	000000-00-0	40



Library Search Compound Report

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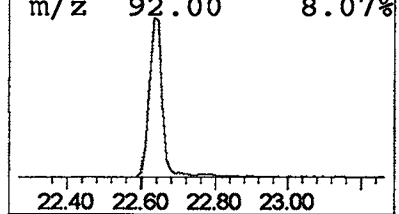
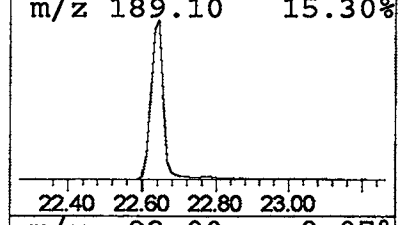
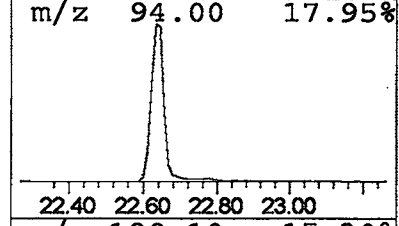
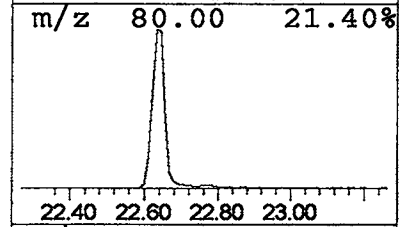
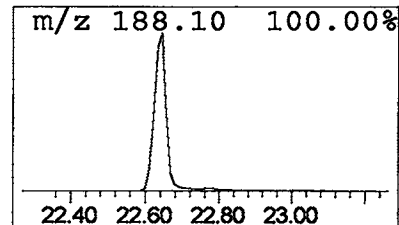
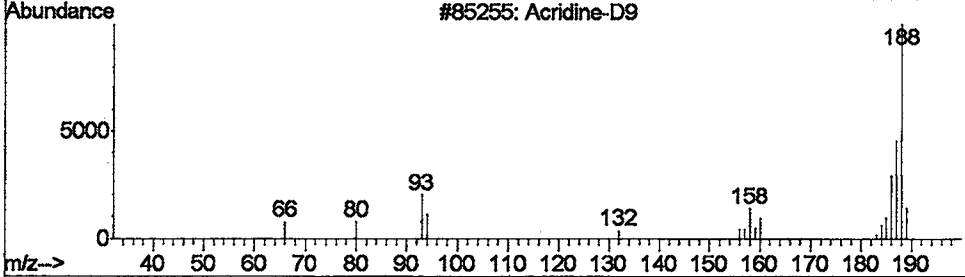
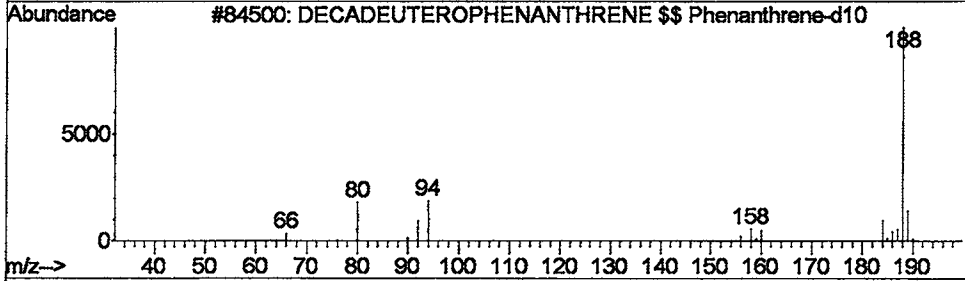
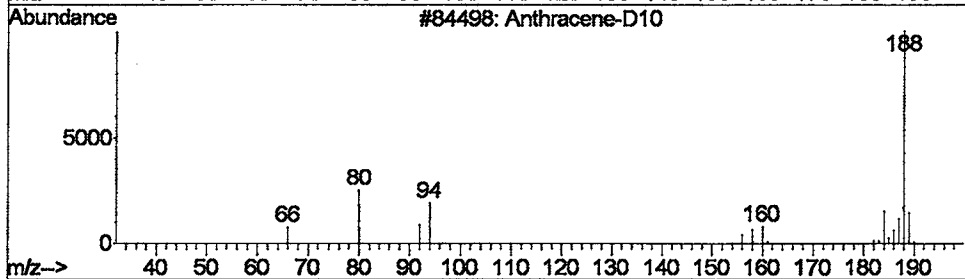
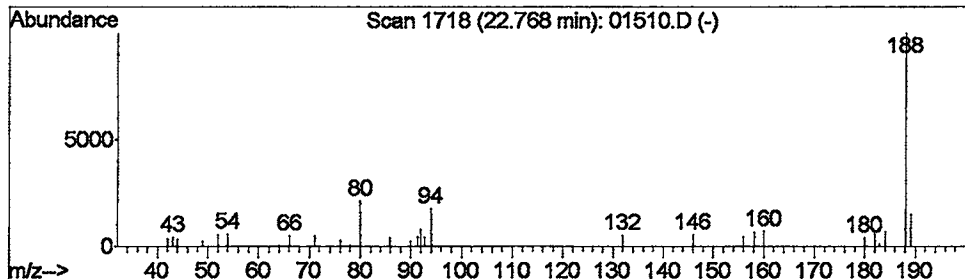
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Quant Method : C:\MSDCHEM\1...\HP022102.M (RTE Integrator)
 Title : 508 scan method
 Library : C:\DATABASE\WILEY7N.L

 Peak Number 11 Anthracene-D10 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.77	0.69 ug/L	529173	Phenanthrene-d10	22.64

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Anthracene-D10	188	C14D10	000000-00-0	90
2			DECADEUTEROPHENANTHRENE \$\$ Phena...	188	C14D10	001517-22-2	83
3			Acridine-D9	188	C13D9N	000000-00-0	64
4			Phenanthrene-d10	188	C14D10	001517-22-2	64



Library Search Compound Report

Data File : C:\MSDCHEM\1\5973N\02FEB21\01510.D
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Sample : 508 scan
Misc : February 21, 2002
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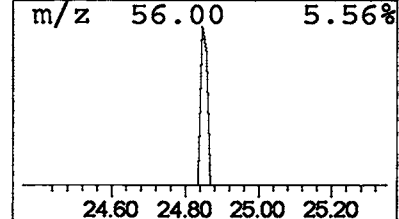
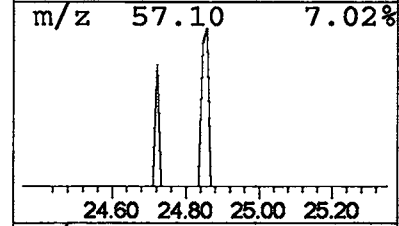
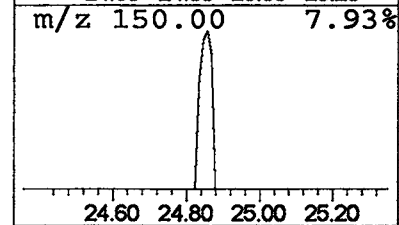
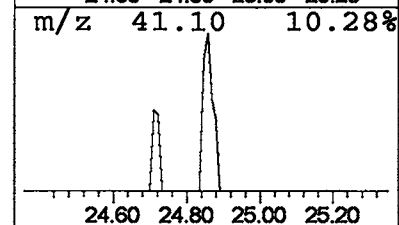
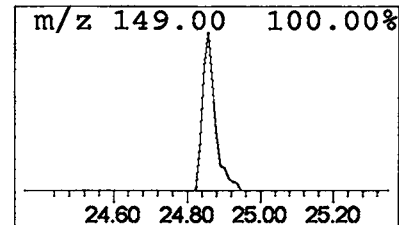
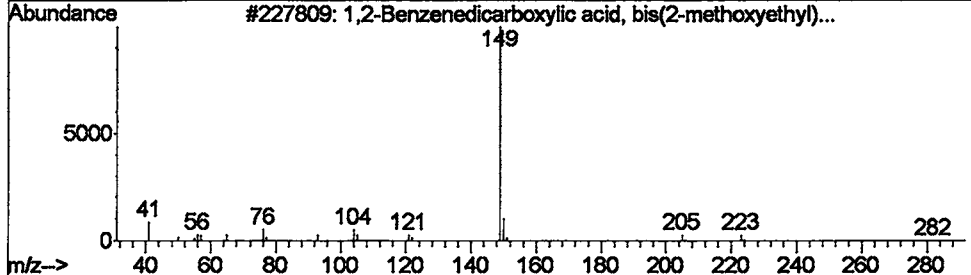
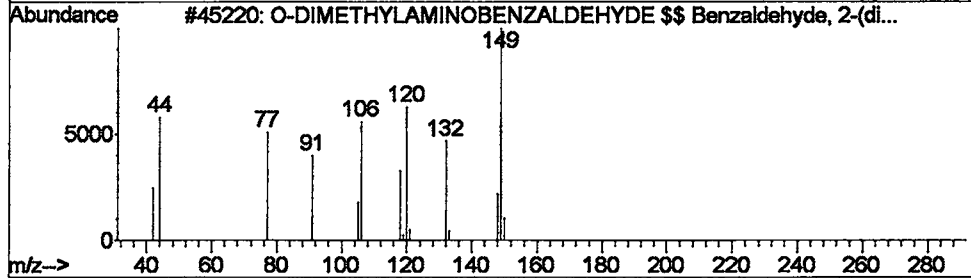
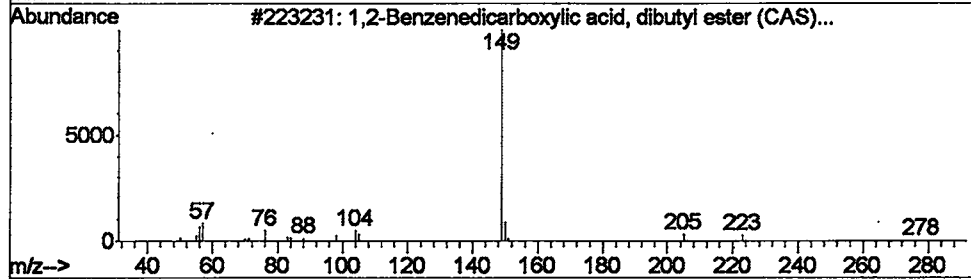
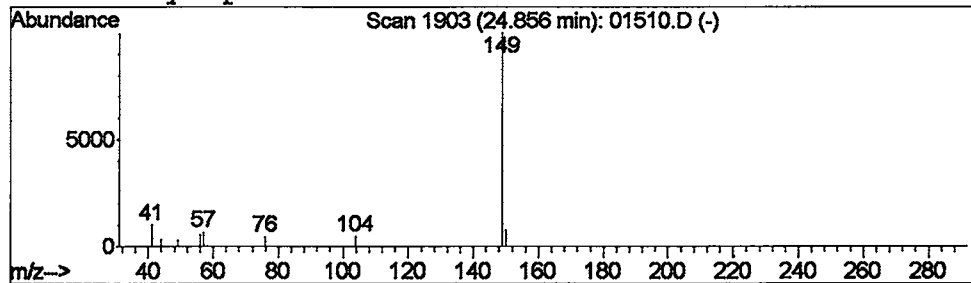
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Library : C:\DATABASE\WILEY7N.L

Peak Number 12 1,2-Benzenedicarboxylic aci... Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.86	0.04 ug/L	33952	Phenanthrene-d10	22.64

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1,2-Benzenedicarboxylic acid, di...	278	C16H22O4	000084-74-2	64
2			O-DIMETHYLAMINOBENZALDEHYDE \$\$ B...	149	C9H11NO	000579-72-6	9
3			1,2-Benzenedicarboxylic acid, bi...	282	C14H18O6	000117-82-8	9
4			Dibutyl phthalate	278	C16H22O4	000084-74-2	9



Library Search Compound Report

Data File : C:\MSDCHEM\1\5973N\02FEB21\01510.D
 Acq On : 21 Feb 2002 2:17 pm
 Sample : 508 scan
 Misc : February 21, 2002
 MS Integration Params: LSCINT.P

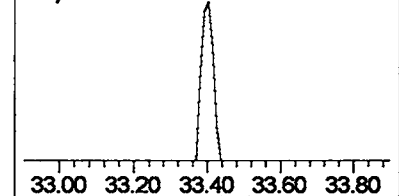
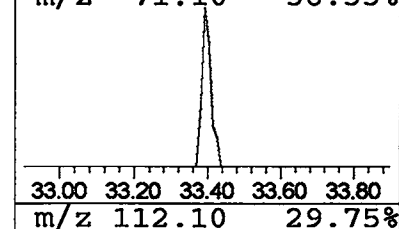
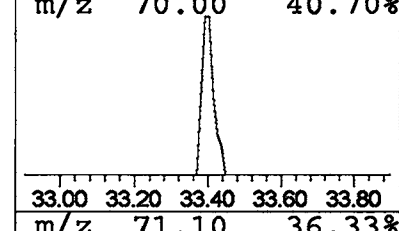
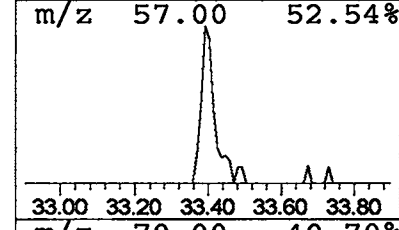
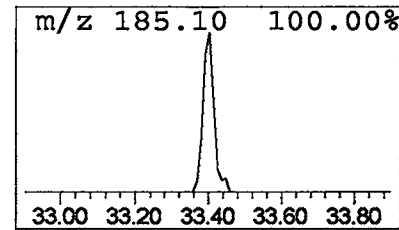
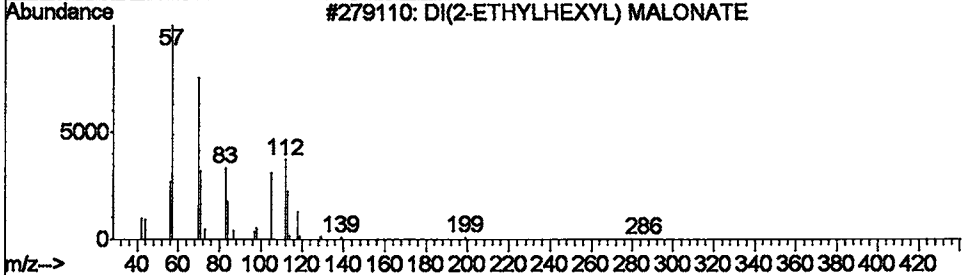
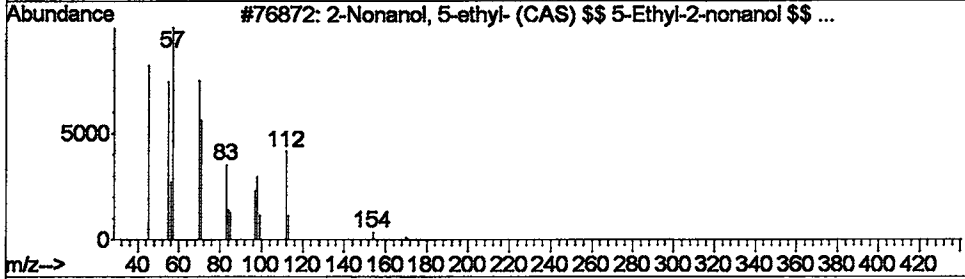
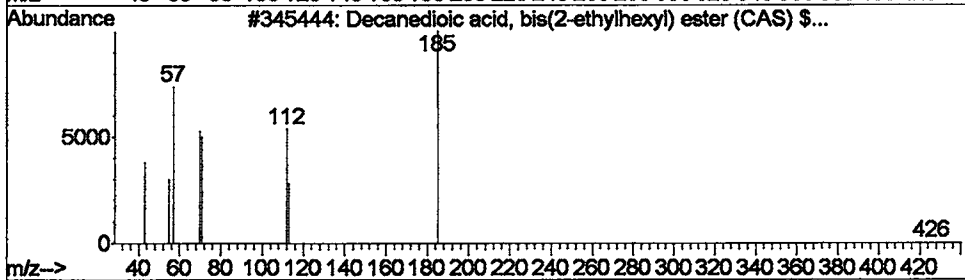
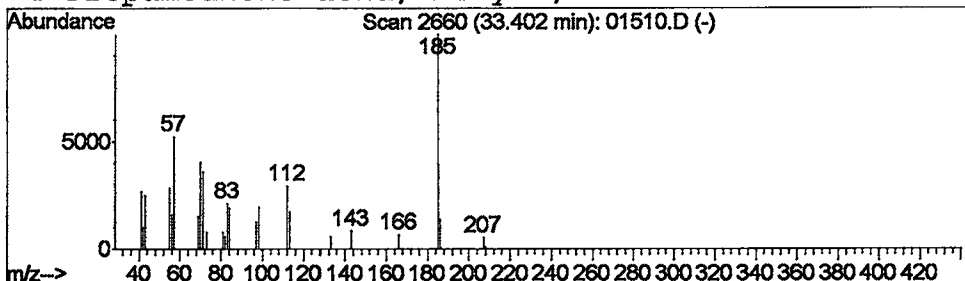
Vial: 5
 Operator:
 Inst : Instrumen
 Multiplr: 1.00

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

 Peak Number 13 Decanedioic acid, bis(2-eth... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
33.40	0.19 ug/L	100968	Perylene-d12	34.41

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Decanedioic acid, bis(2-ethylhex...	426	C26H50O4	000122-62-3	50
2		2-Nonanol, 5-ethyl- (CAS) \$\$ 5-E...	172	C11H24O	000103-08-2	37
3		DI(2-ETHYLHEXYL) MALONATE	328	C19H36O4	000000-00-0	35
4		Propanedioic acid, hexyl-, dieth...	244	C13H24O4	005398-10-7	35



Tentatively Identified Compound (LSC) summary

Operator ID: Date Acquired: 21 Feb 2002 2:17 pm
 Data File: C:\MSDCHEM\1\5973N\02FEB21\01510.D
 Name: 508 scan
 Misc: February 21, 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
Dichloroacetonitr...	3.50	0.1	ug/L	42630	1	9.72	9133220	20.0
Butanoic acid, me...	3.62	0.2	ug/L	70261	1	9.72	9133220	20.0
ISOBUTYRALDEHYDE,...	3.93	0.1	ug/L	37728	1	9.72	9133220	20.0
1,3,5-Cycloheptat...	4.38	0.1	ug/L	32697	1	9.72	9133220	20.0
3-(CYCLOHEX-3'-EN...	5.95	0.5	ug/L	240299	1	9.72	9133220	20.0
2-Pentanone, 4-hy...	6.06	0.5	ug/L	227971	1	9.72	9133220	20.0
2-Propanol, 1-(2-...	9.88	0.2	ug/L	77760	1	9.72	9133220	20.0
2,5-Cyclohexadien...	13.38	0.0	ug/L	32628	2	13.20	13611400	20.0
3-Buten-2-ol, 2-m...	19.98	0.1	ug/L	51884	3	18.34	14701100	20.0
Phenanthrene, 9-m...	22.28	0.1	ug/L	98114	4	22.64	15439200	20.0
Anthracene-D10	22.77	0.7	ug/L	529173	4	22.64	15439200	20.0
1,2-Benzenedicarb...	24.86	0.0	ug/L	33952	4	22.64	15439200	20.0
Decanedioic acid,...	33.40	0.2	ug/L	100968	6	34.41	10563200	20.0