

User: Vinci, David L
Westchester Dept. of Labs & Research
Date: 02/27/2002 Time: 09:07:09

Sample: AE01142
Analysis: \$GCMS GCMS Scan For Unknowns
Units: ug
Started: 2/27/02 09:06 Ended: 2/27/02 09:06 Analyst: DLV
Page: 1

	Component Name	Viol	Result	Qualifier	MDL
1	Other unknown compounds		.		
2	Surr_2,4-DDD		98		5.0

Comments for Sample AE01142 - Analysis \$GCMS

Westchester Dept. of Labs & Research

User: Vinci, David L

Date: 02-27-2002 Time: 09:07:12

The GCMS scan, from 35 to 550 amu, does not reveal the presence of unusual compounds. The recoveries for 16 of the 43 compounds in the LCS Standard were above their acceptance ranges, while 3 of the 43 compounds were below their acceptance ranges.

Data File : C:\MSDCHEM\1\5973N\02FEB25\01142.D

Vial: 8

Acq On : 25 Feb 2002 5:01 pm

Operator:

Sample : 508 scan

Inst : Instrumen

Misc : February 25, 2002

Multiplr: 1.00

MS Integration Params: SPLIT.P

Quant Time: Feb 26 08:28:42 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	1536734	20.00	ug/L	0.00
10) Naphthalene-d8	13.20	136	6522782	20.00	ug/L	0.00
17) Acenaphthene-d10	18.33	164	3186051	20.00	ug/L	0.00
29) Phenanthrene-d10	22.63	188	5117214	20.00	ug/L	-0.01
38) Chrysene-d12	30.49	240	4119675	20.00	ug/L	-0.03
44) Perylene-d12	34.41	264	2920714	20.00	ug/L	-0.04

System Monitoring Compounds

37) 2,4-DDD	27.72	235	335184	4.91	ug/L	0.00
Spiked Amount	5.000		Recovery	=	98.20%	

Target Compounds

						Qvalue
20) Dimethylphthalate	18.34	163	513990	2.66	ug/L	# 58
21) 2,6-Dinitrotoluene	18.33	165	407448	9.13	ug/L	# 27

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

01142.D HP022102.M Tue Feb 26 08:28:43 2002

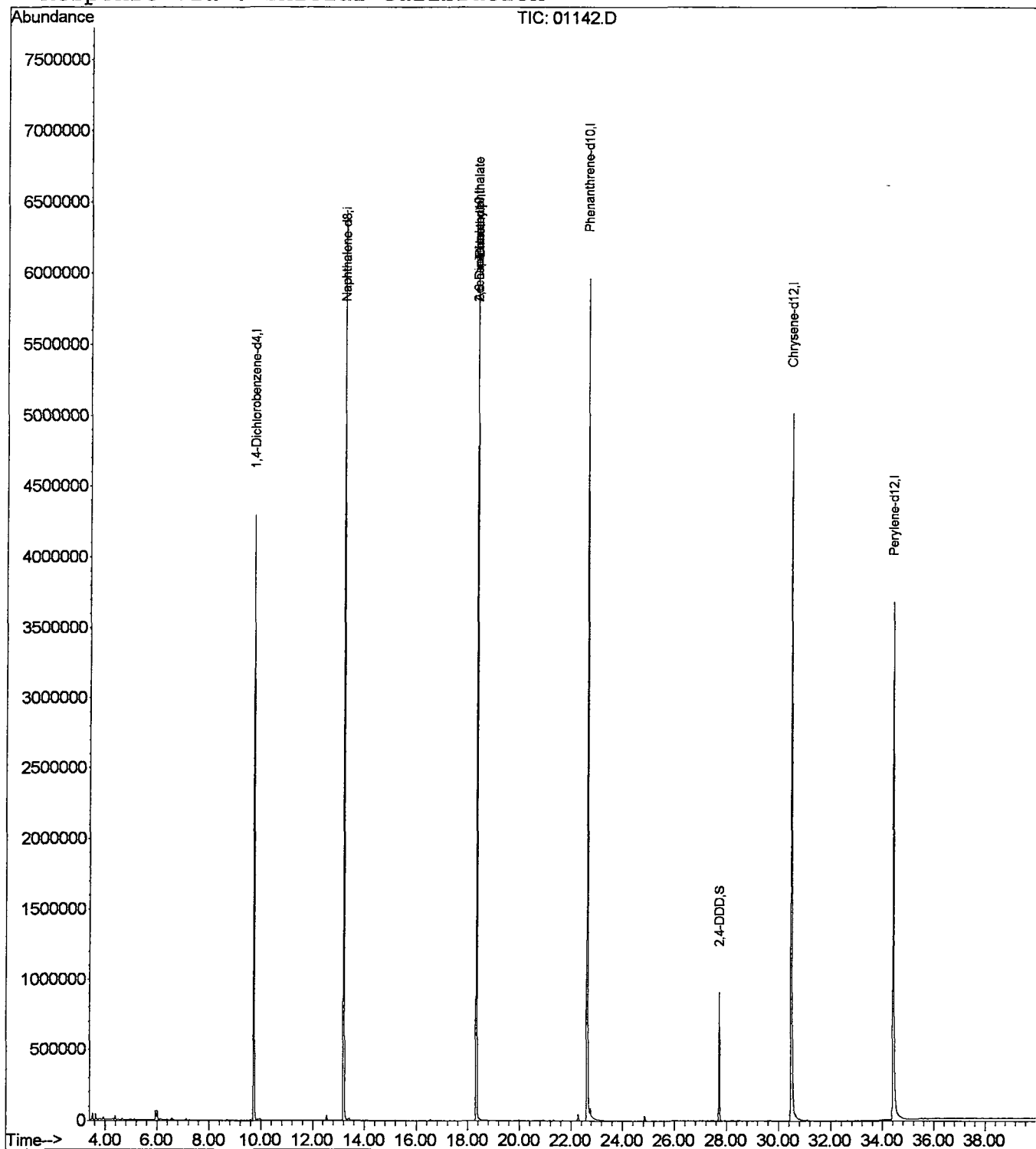
Page 1

Data File : C:\MSDCHEM\1\5973N\02FEB25\01142.D
 Acq On : 25 Feb 2002 5:01 pm
 Sample : 508 scan
 Misc : February 25, 2002
 MS Integration Params: SPLIT.P
 Quant Time: Feb 26 8:28 2002

Vial: 8
 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



Data File : C:\MSDCHEM\1\5973N\02FEB25\01142.D

Vial: 8

Acq On : 25 Feb 2002 5:01 pm

Operator:

Sample : 508 scan

Inst : Instrumen

Misc : February 25, 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 >

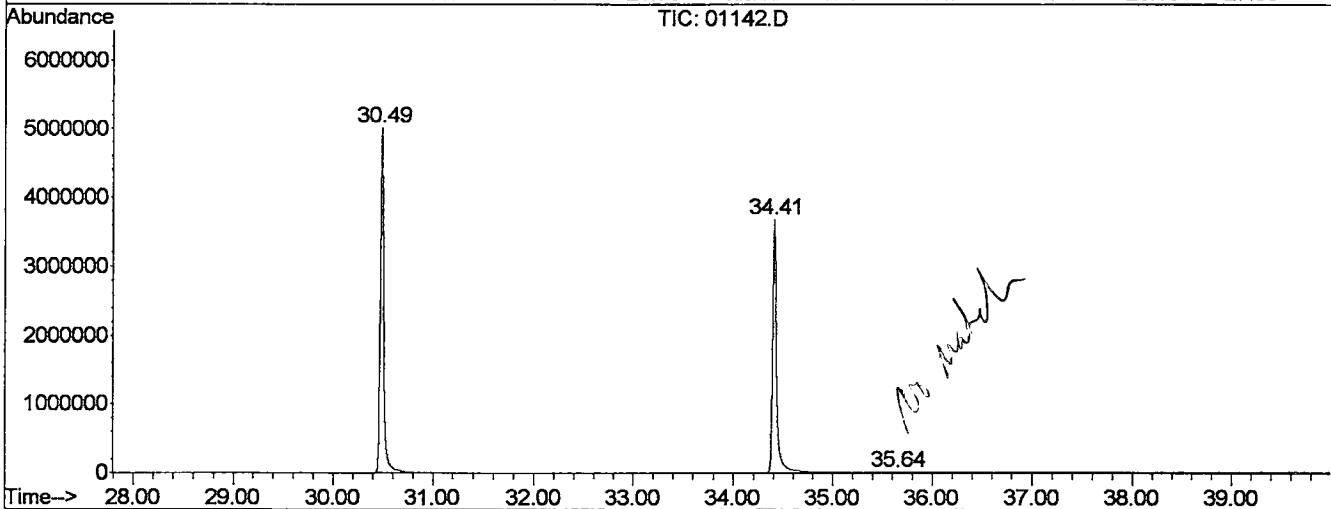
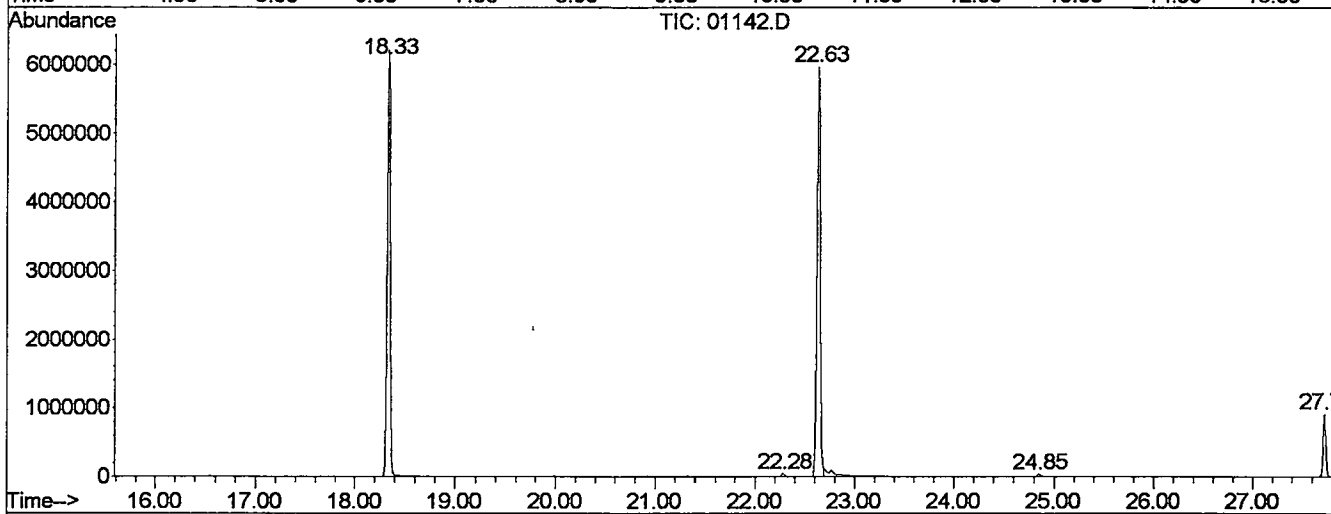
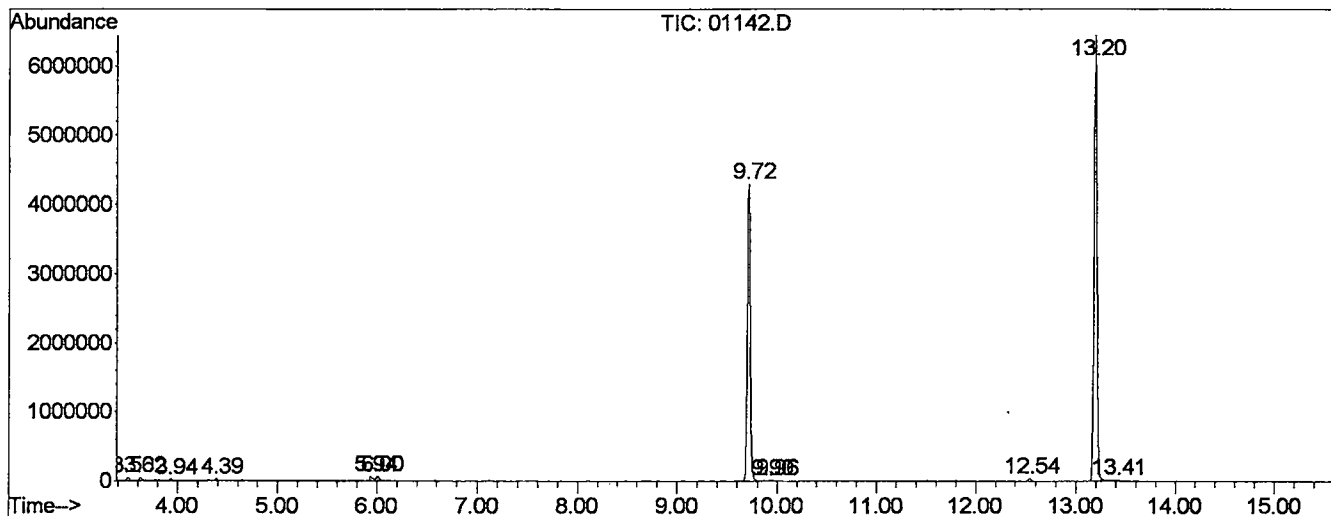
Peak separation: 5

Signal : TIC

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.498	9	11	15	rVB	44089	73427	0.55%	0.102%
2	3.622	19	22	25	rBV	40295	60193	0.45%	0.084%
3	3.938	45	50	53	rVB	17418	32748	0.25%	0.046%
4	4.390	87	90	93	rBV	27728	48469	0.36%	0.067%
5	5.936	223	227	231	rBV	65358	137872	1.03%	0.192%
6	6.004	231	233	241	rVV	66118	138635	1.04%	0.193%
7	9.718	557	562	567	rBV	4294168	8534006	63.98%	11.871%
8	9.899	573	578	581	rBV2	7637	28692	0.22%	0.040%
9	9.955	581	583	591	rVB2	13211	26666	0.20%	0.037%
10	12.540	809	812	817	rBV	36163	62986	0.47%	0.088%
11	13.195	865	870	875	rBV	6437133	12350667	92.60%	17.180%
12	13.410	883	889	895	rVB2	15704	50520	0.38%	0.070%
13	18.332	1319	1325	1331	rBV	6211423	13006835	97.52%	18.093%
14	22.283	1671	1675	1683	rVB	44423	89300	0.67%	0.124%
15	22.633	1701	1706	1711	rBV2	5963038	13338101	100.00%	18.554%
16	24.845	1897	1902	1915	rVB	32035	79361	0.59%	0.110%
17	27.724	2151	2157	2163	rBV	909946	1639862	12.29%	2.281%
18	30.490	2397	2402	2435	rBV2	5010338	12189300	91.39%	16.956%
19	34.407	2743	2749	2787	rBV	3674577	9977560	74.80%	13.879%
20	35.638	2855	2858	2865	rBV6	6460	24017	0.18%	0.033%

Sum of corrected areas: 71889217

File : C:\MSDCHEM\1\5973N\02FEB25\01142.D
 Operator :
 Acquired : 25 Feb 2002 5:01 pm using AcqMethod HP020702
 Instrument : Instrumen
 Sample Name: 508 scan
 Misc Info : February 25, 2002
 Vial Number: 8
 Quant File :HP022102.RES (RTE Integrator)



Data File : C:\MSDCHEM\1\5973N\02FEB25\01142.D
 Acq On : 25 Feb 2002 5:01 pm
 Sample : 508 scan
 Misc : February 25, 2002
 MS Integration Params: LSCINT.P

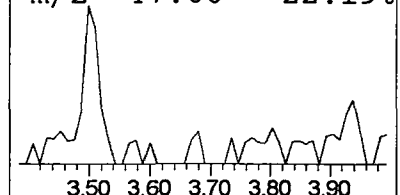
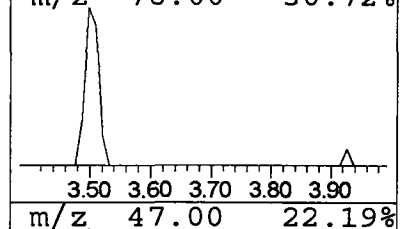
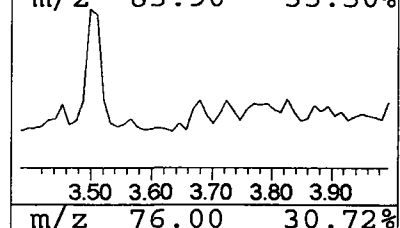
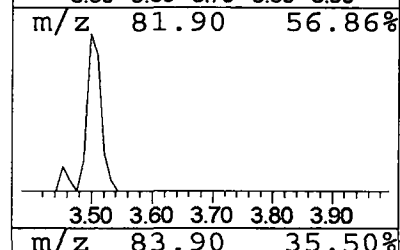
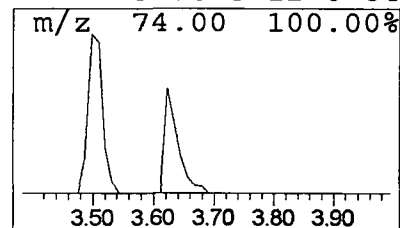
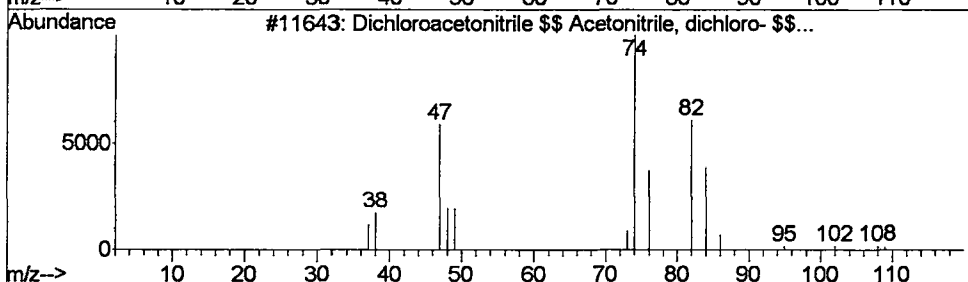
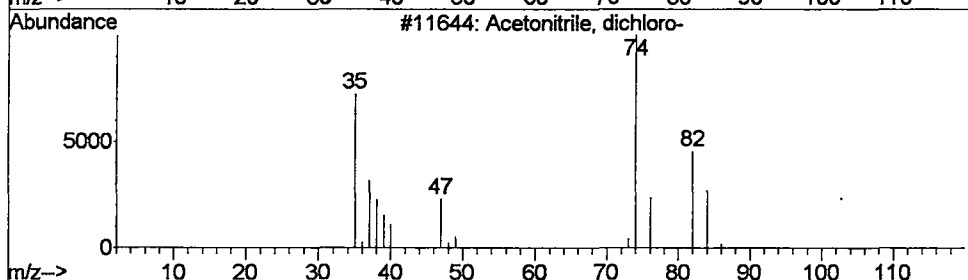
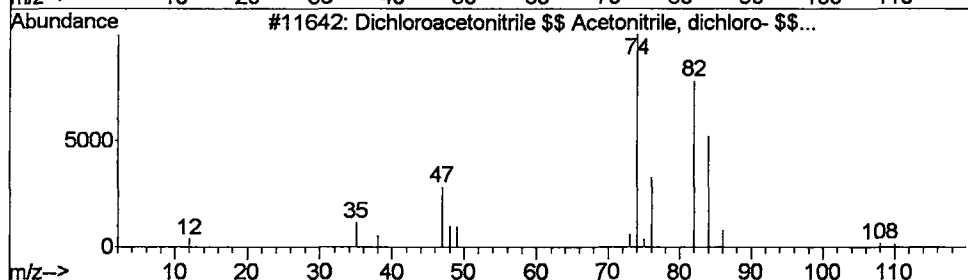
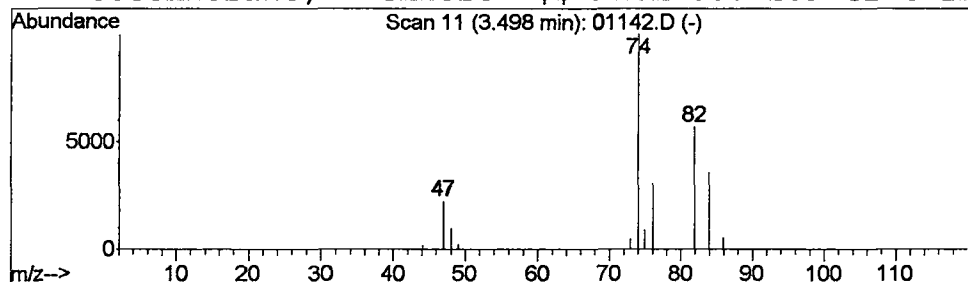
Vial: 8
 Operator:
 Inst : Instrumen
 Multiplr: 1.00

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 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.50	0.17 ug/L	73427	1,4-Dichlorobenzene-d4	9.72

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Dichloroacetonitrile \$\$ Acetonit...	109	C2HCl2N	003018-12-0	83
2		Acetonitrile, dichloro-	109	C2HCl2N	003018-12-0	74
3		Dichloroacetonitrile \$\$ Acetonit...	109	C2HCl2N	003018-12-0	74
4		Acetonitrile, dichloro- \$\$ Dichl...	109	C2HCl2N	003018-12-0	64



Data File : C:\MSDCHEM\1\5973N\02FEB25\01142.D
 Acq On : 25 Feb 2002 5:01 pm
 Sample : 508 scan
 Misc : February 25, 2002
 MS Integration Params: LSCINT.P

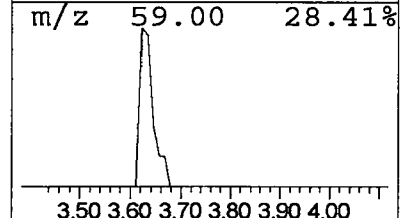
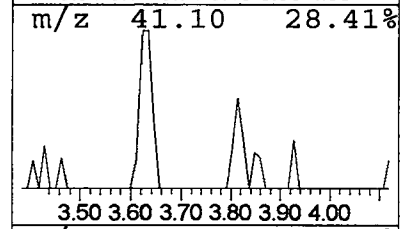
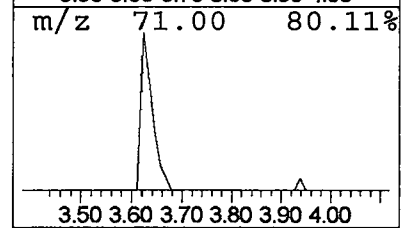
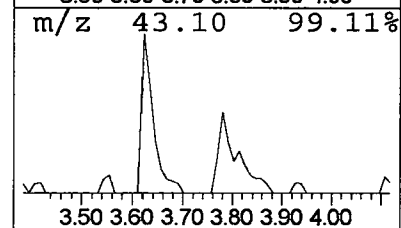
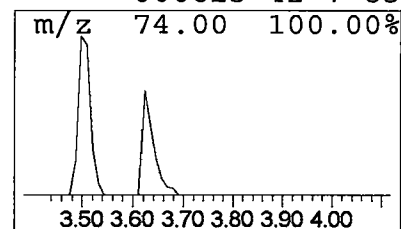
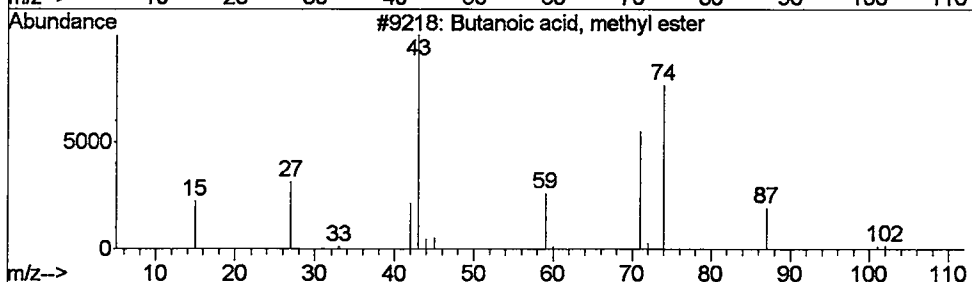
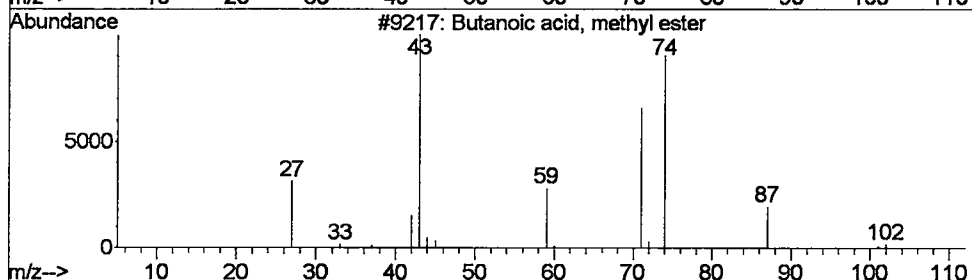
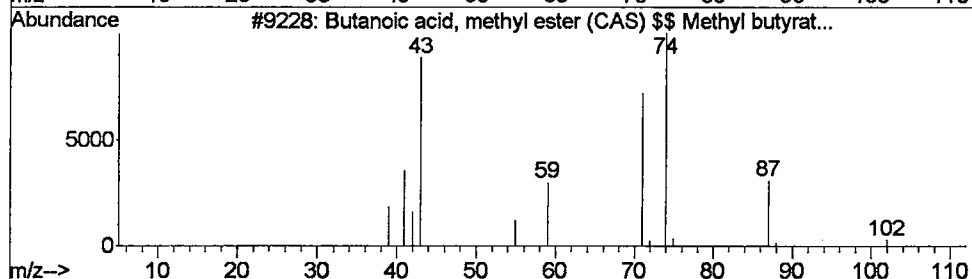
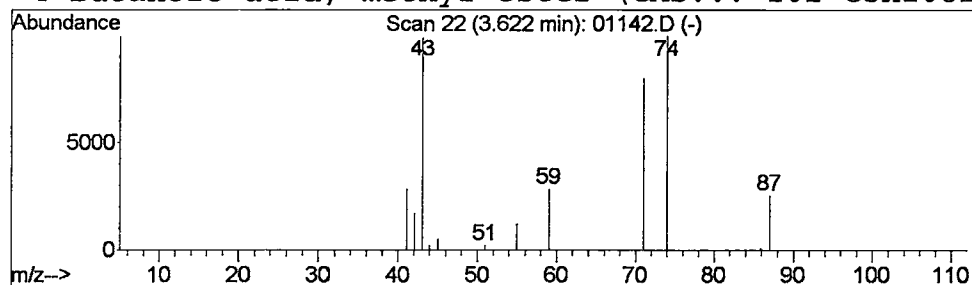
Vial: 8
 Operator:
 Inst : Instrumen
 Multiplr: 1.00

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	0.14 ug/L	60193	1,4-Dichlorobenzene-d4	9.72

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	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



Data File : C:\MSDCHEM\1\5973N\02FEB25\01142.D
 Acq On : 25 Feb 2002 5:01 pm
 Sample : 508 scan
 Misc : February 25, 2002
 MS Integration Params: LSCINT.P

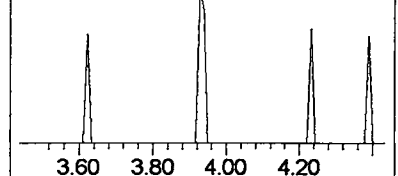
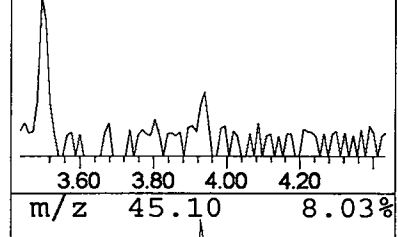
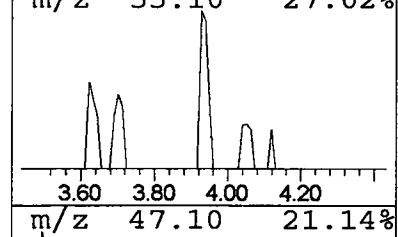
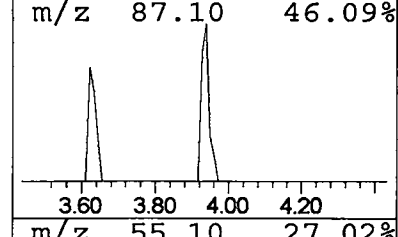
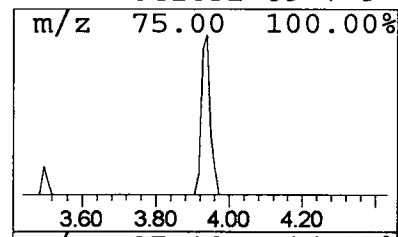
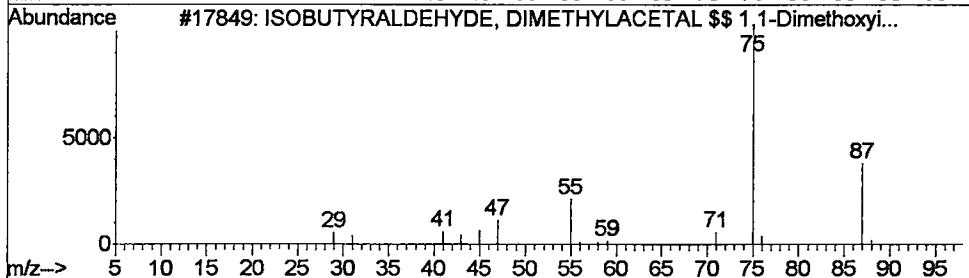
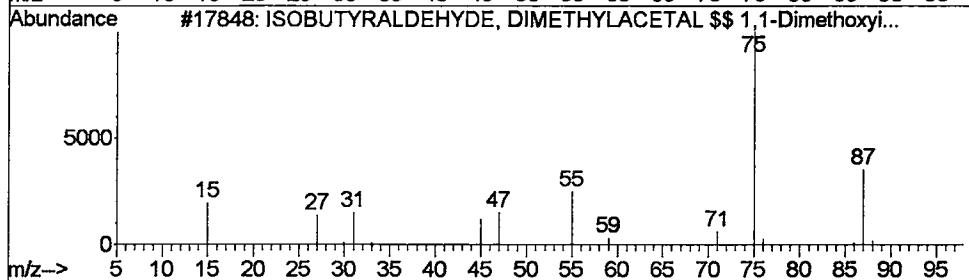
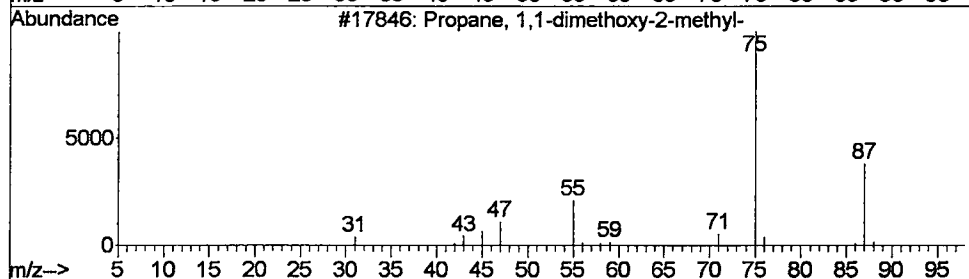
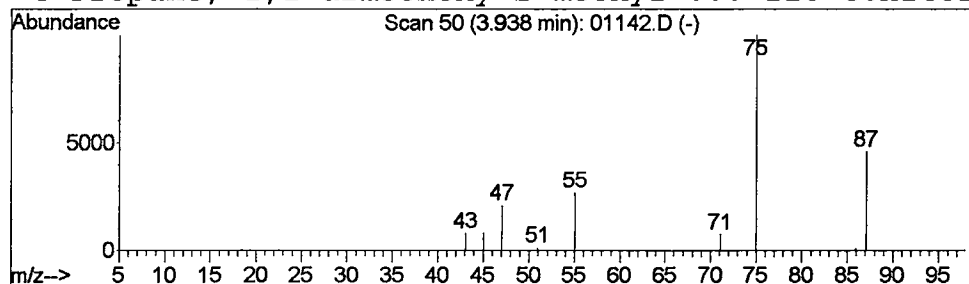
Vial: 8
 Operator:
 Inst : Instrumen
 Multiplr: 1.00

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 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.94	0.08 ug/L	32748	1,4-Dichlorobenzene-d4	9.72

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Propane, 1,1-dimethoxy-2-methyl-	118	C6H14O2	041632-89-7	78
2			ISOBUTYRALDEHYDE, DIMETHYLACETAL...	118	C6H14O2	000000-00-0	64
3			ISOBUTYRALDEHYDE, DIMETHYLACETAL...	118	C6H14O2	000000-00-0	43
4			Propane, 1,1-dimethoxy-2-methyl-...	118	C6H14O2	041632-89-7	9



Data File : C:\MSDCHEM\1\5973N\02FEB25\01142.D
 Acq On : 25 Feb 2002 5:01 pm
 Sample : 508 scan
 Misc : February 25, 2002
 MS Integration Params: LSCINT.P

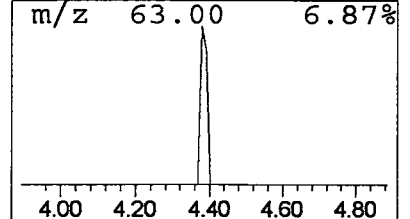
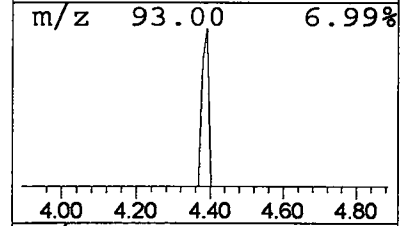
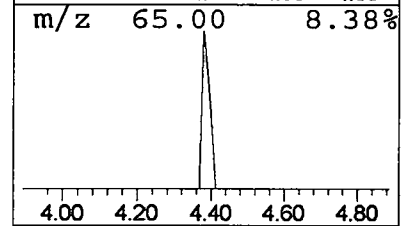
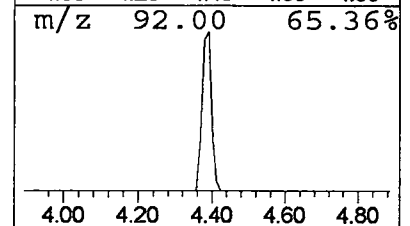
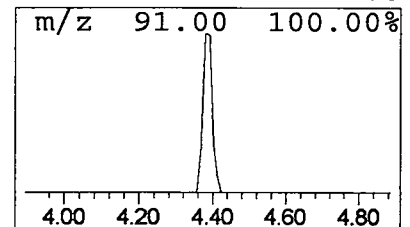
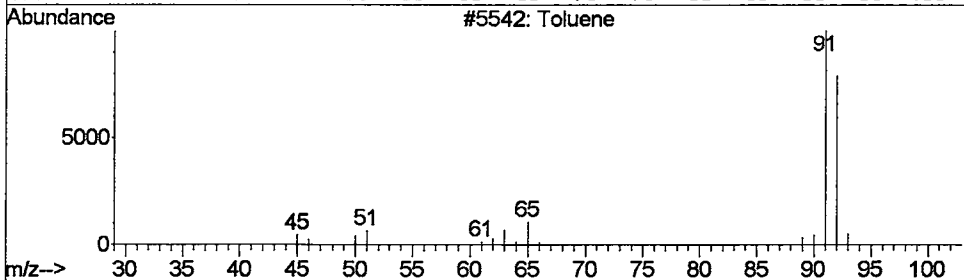
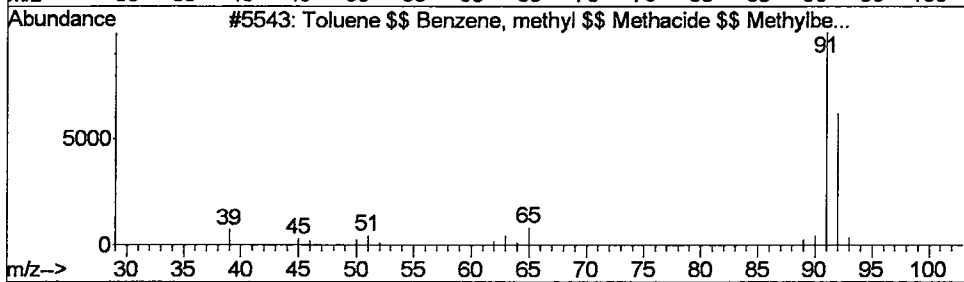
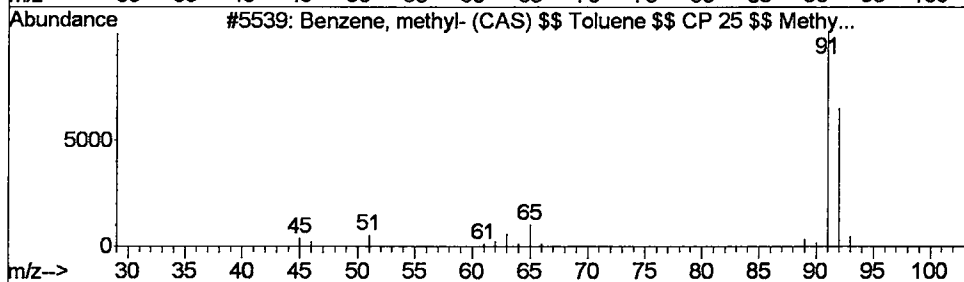
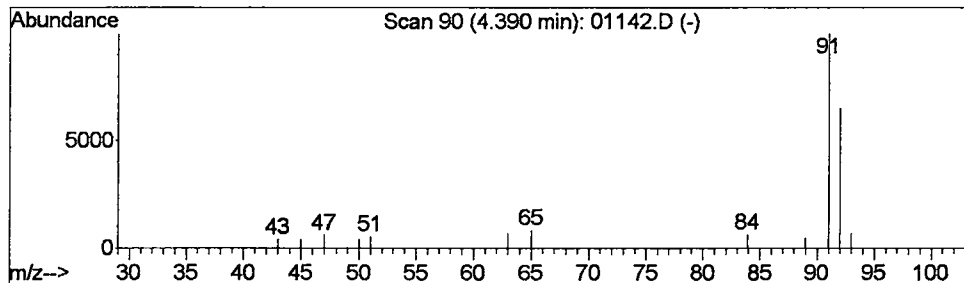
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 Operator:
 Inst : Instrumen
 Multiplr: 1.00

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

 Peak Number 4 Benzene, methyl- (CAS) \$\$ T... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.39	0.11 ug/L	48469	1,4-Dichlorobenzene-d4	9.72

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, methyl- (CAS) \$\$ Toluen...	92	C7H8	000108-88-3	86
2		Toluene \$\$ Benzene, methyl \$\$ Me...	92	C7H8	000108-88-3	86
3		Toluene	92	C7H8	000108-88-3	86
4		Toluene	92	C7H8	000108-88-3	80



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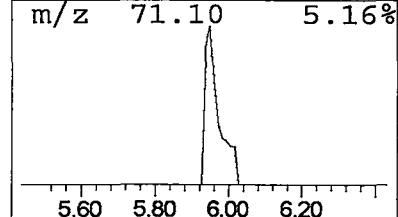
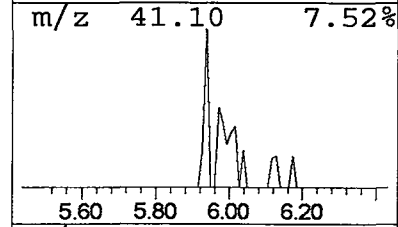
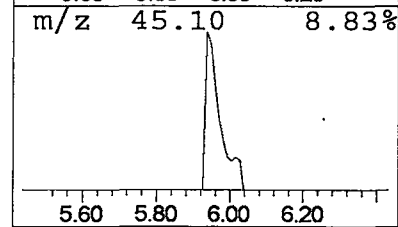
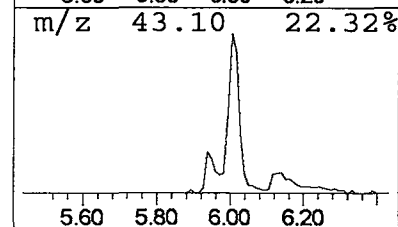
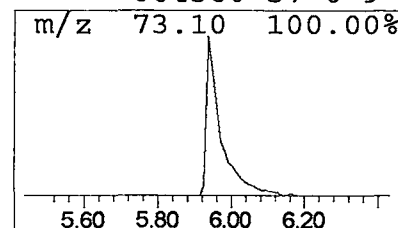
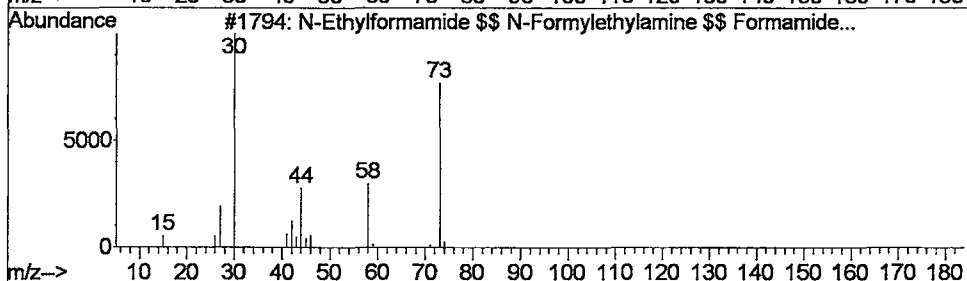
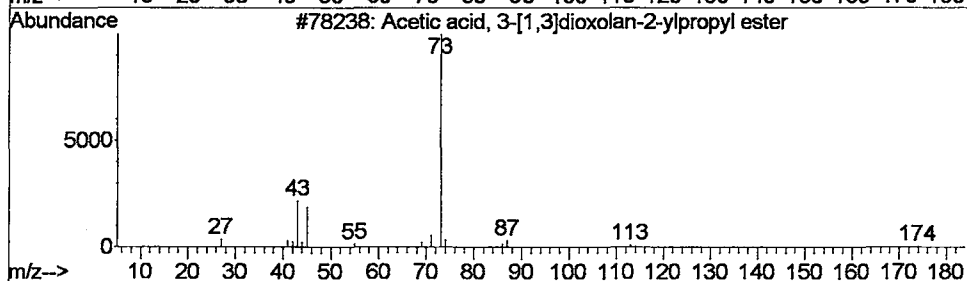
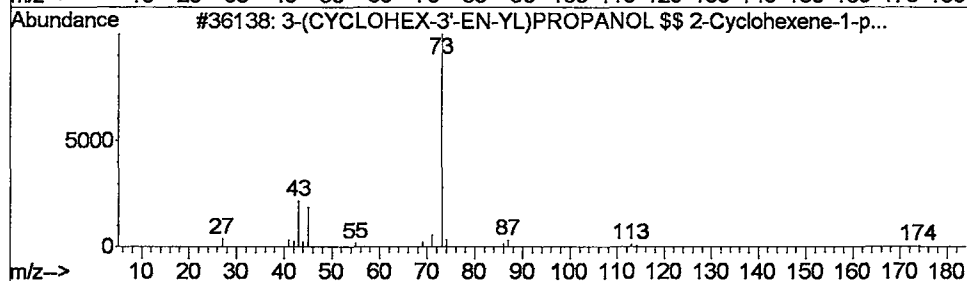
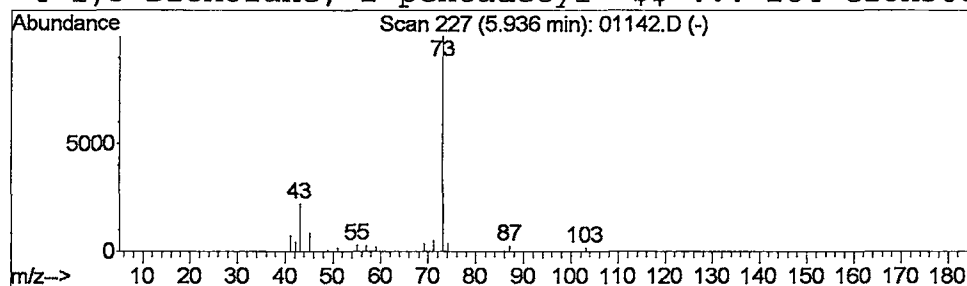
Vial: 8
 Operator:
 Inst : Instrumen
 Multiplr: 1.00

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.
5.94	0.32 ug/L	137872	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	38
2		Acetic acid, 3-[1,3]dioxolan-2-y...	174	C8H14O4	000000-00-0	38
3		N-Ethylformamide \$\$ N-Formylethy...	73	C3H7NO	000627-45-2	9
4		1,3-Dioxolane, 2-pentadecyl- \$\$...	284	C18H36O2	004360-57-0	9



Data File : C:\MSDCHEM\1\5973N\02FEB25\01142.D
 Acq On : 25 Feb 2002 5:01 pm
 Sample : 508 scan
 Misc : February 25, 2002
 MS Integration Params: LSCINT.P

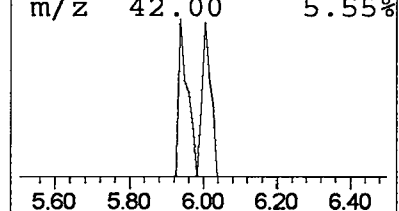
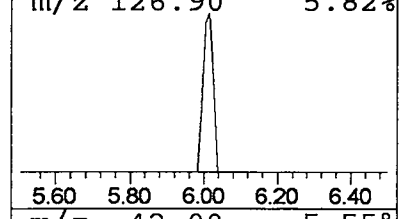
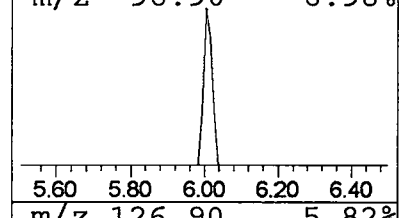
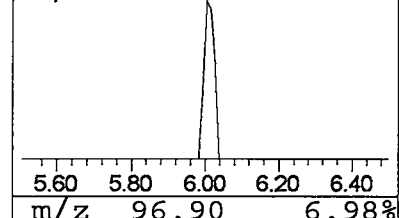
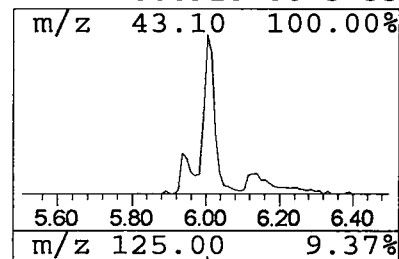
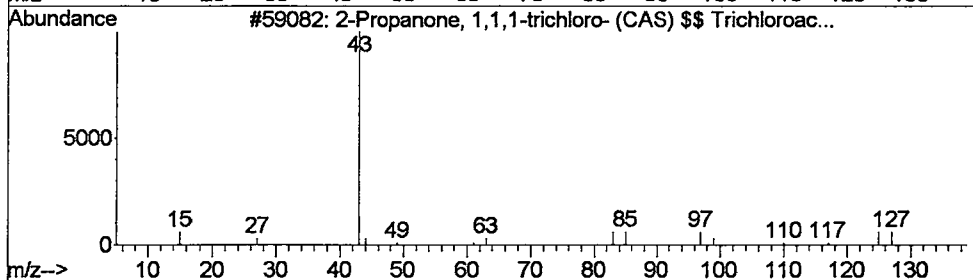
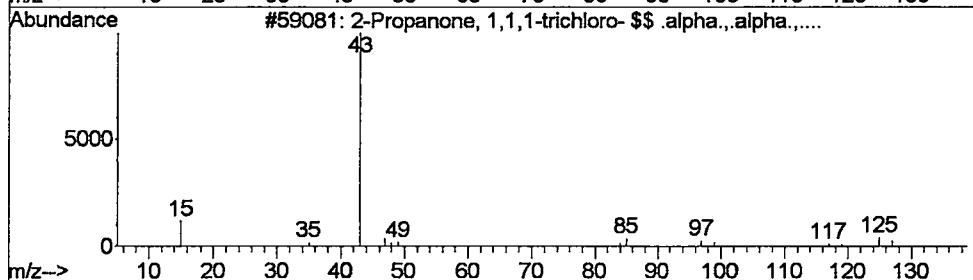
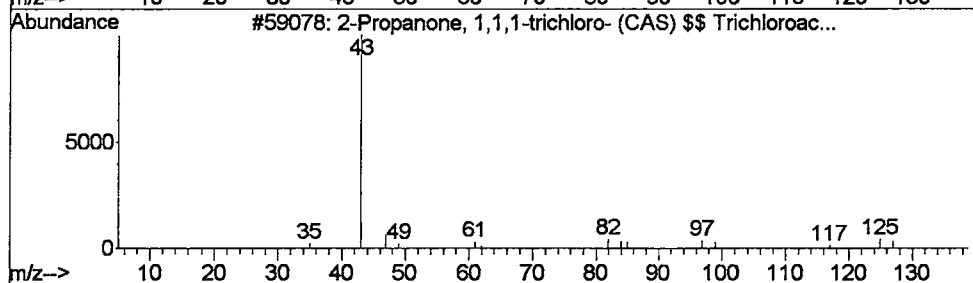
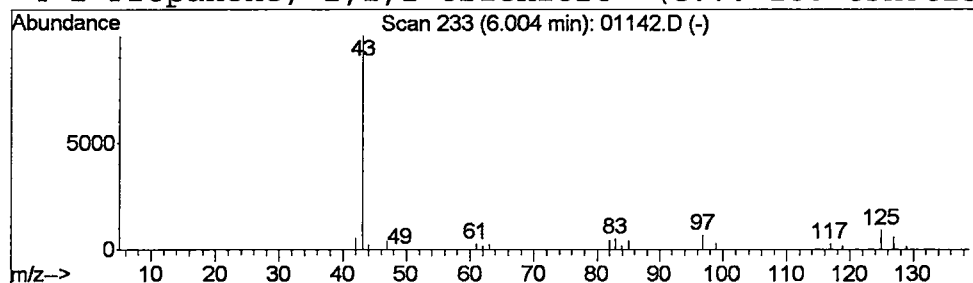
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 Inst : Instrumen
 Multiplr: 1.00

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

 Peak Number 6 2-Propanone, 1,1,1-trichloro... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.00	0.32 ug/L	138635	1,4-Dichlorobenzene-d4	9.72

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Propanone, 1,1,1-trichloro- (C...	160	C3H3Cl3O	000918-00-3	50
2			2-Propanone, 1,1,1-trichloro- \$\$...	160	C3H3Cl3O	000918-00-3	50
3			2-Propanone, 1,1,1-trichloro- (C...	160	C3H3Cl3O	000918-00-3	42
4			2-Propanone, 1,1,1-trichloro- (C...	160	C3H3Cl3O	000918-00-3	33



Data File : C:\MSDCHEM\1\5973N\02FEB25\01142.D
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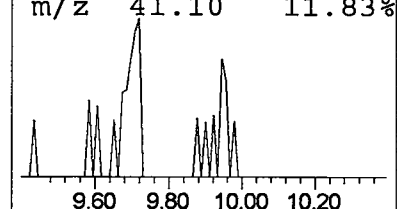
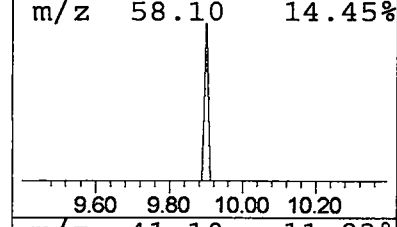
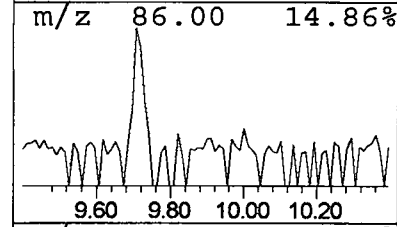
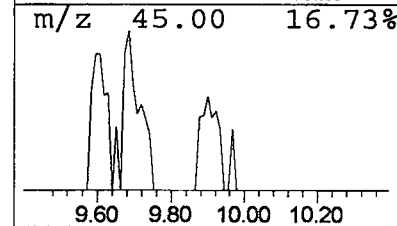
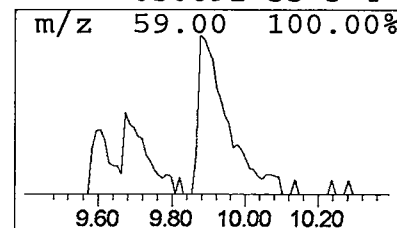
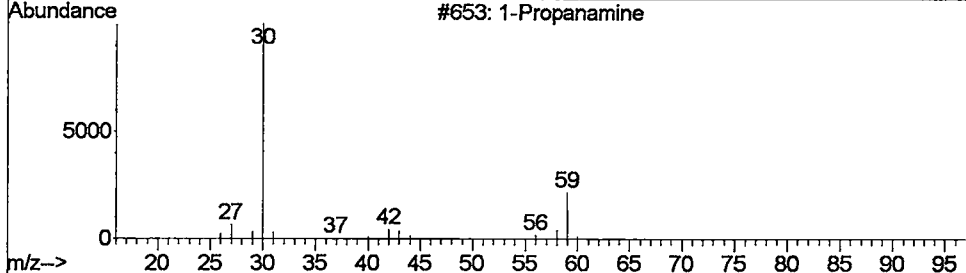
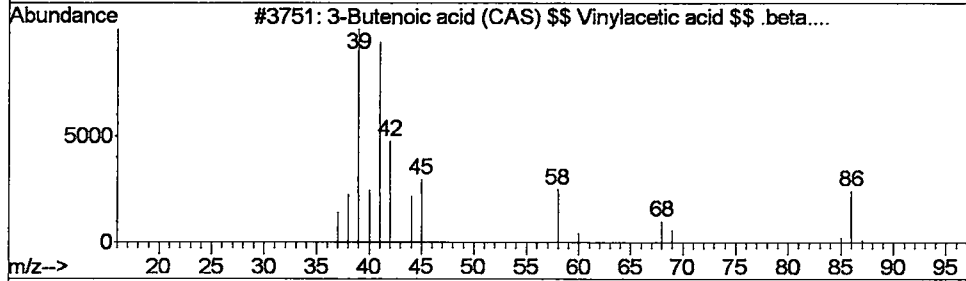
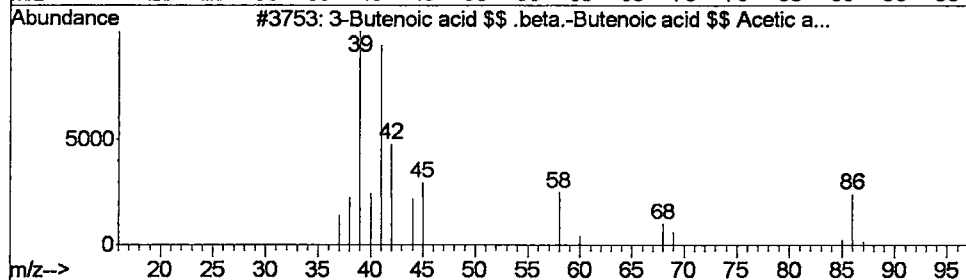
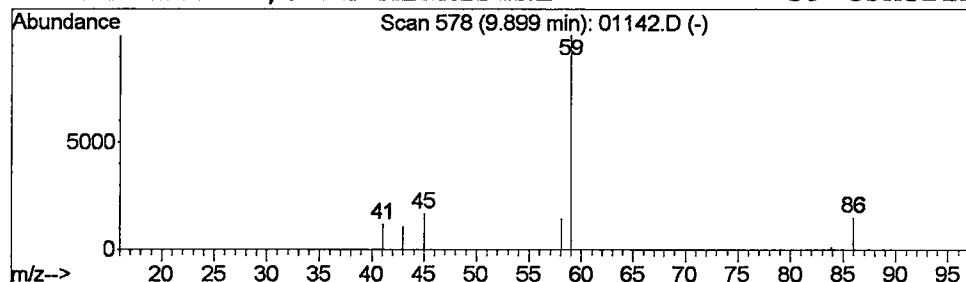
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 Multiplr: 1.00

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

 Peak Number 7 3-Butenoic acid \$\$.beta.-B... Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.90	0.07 ug/L	28692	1,4-Dichlorobenzene-d4	9.72

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			3-Butenoic acid \$\$.beta.-Buteno...	86	C4H6O2	000625-38-7	7
2			3-Butenoic acid (CAS) \$\$ Vinylac...	86	C4H6O2	000625-38-7	7
3			1-Propanamine	59	C3H9N	000107-10-8	5
4			2-METHYL-3,3-D2-AZIRIDINE	59	C3H5D2N	030691-55-5	4



Data File : C:\MSDCHEM\1\5973N\02FEB25\01142.D
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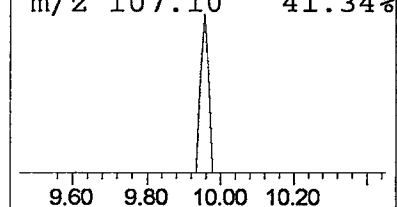
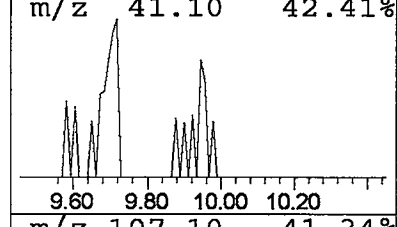
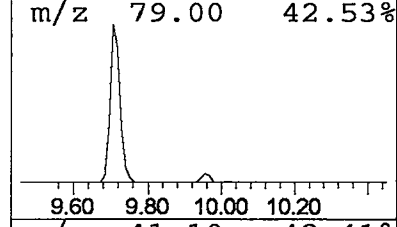
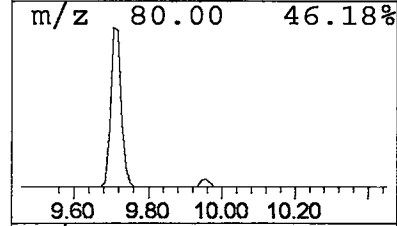
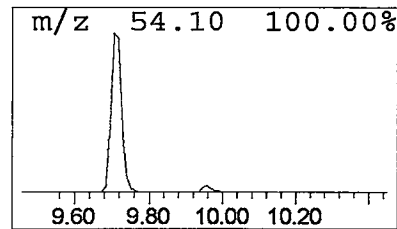
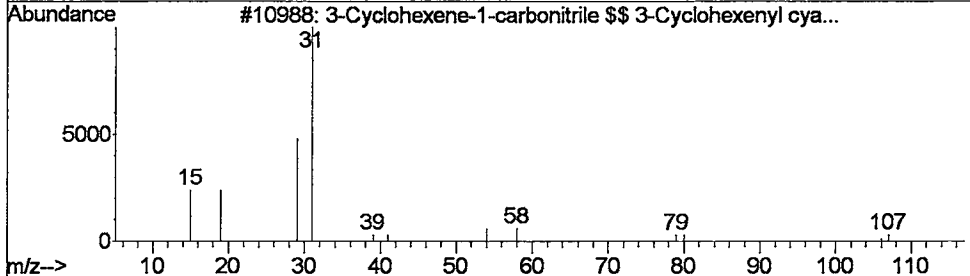
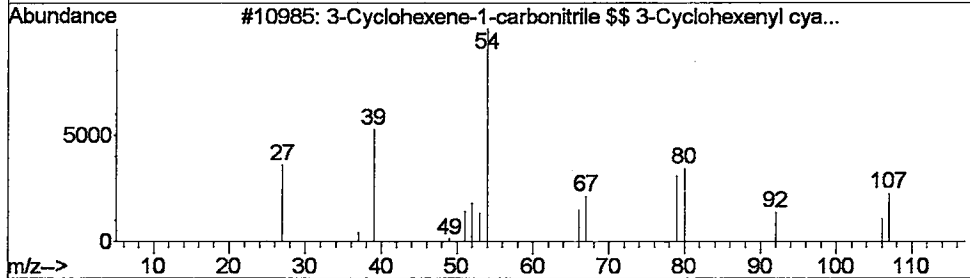
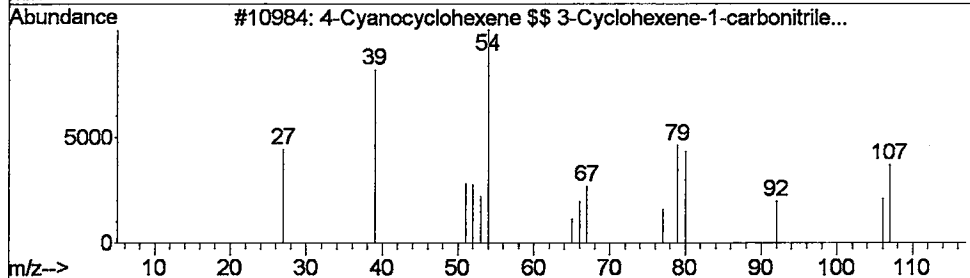
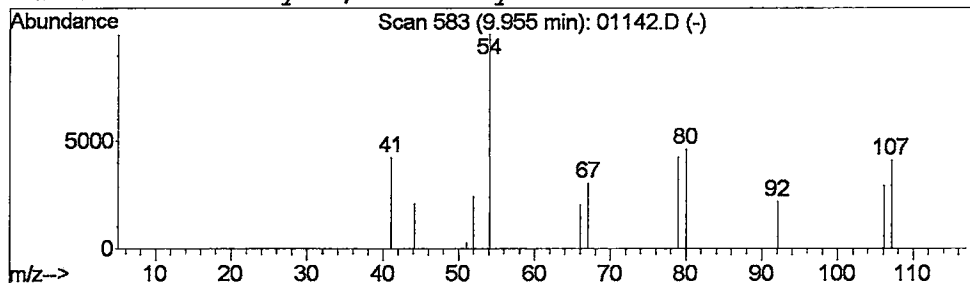
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 Multiplr: 1.00

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

 Peak Number 8 4-Cyanocyclohexene \$\$ 3-Cyc... Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.96	0.06 ug/L	26666	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	91
2		3-Cyclohexene-1-carbonitrile \$\$...	107	C7H9N	000100-45-8	80
3		3-Cyclohexene-1-carbonitrile \$\$...	107	C7H9N	000100-45-8	9
4		2-Penten-4-yne, 2-methyl-	80	C6H8	000000-00-0	9



Data File : C:\MSDCHEM\1\5973N\02FEB25\01142.D
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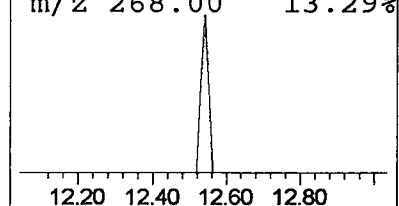
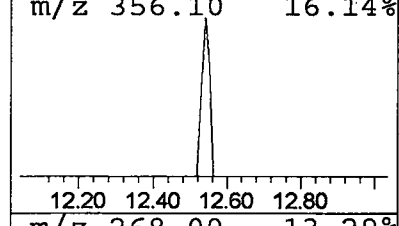
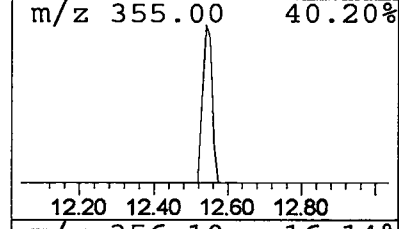
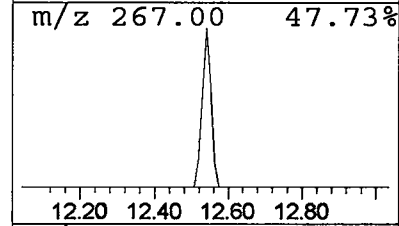
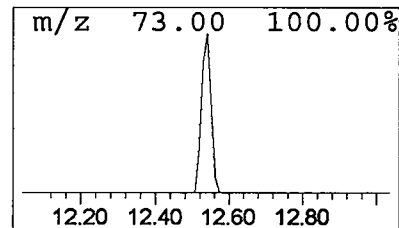
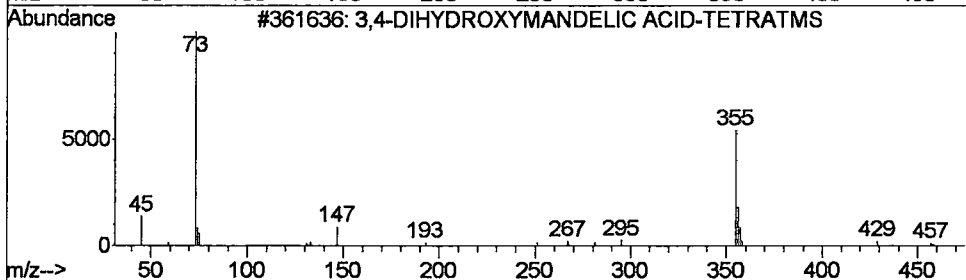
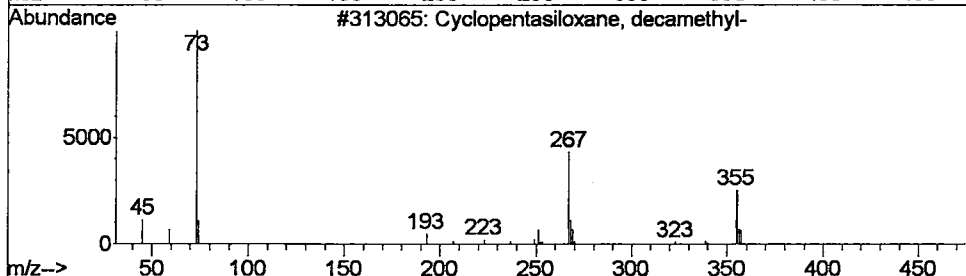
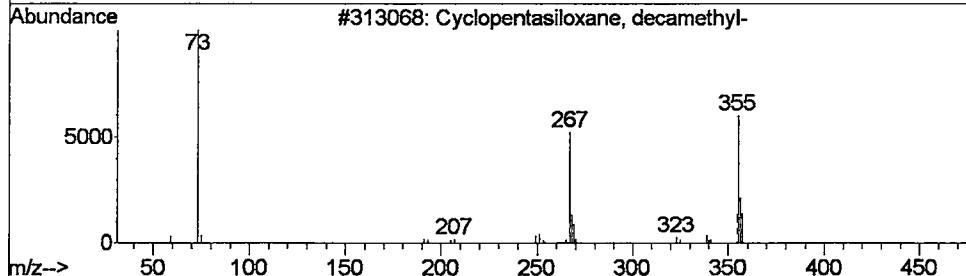
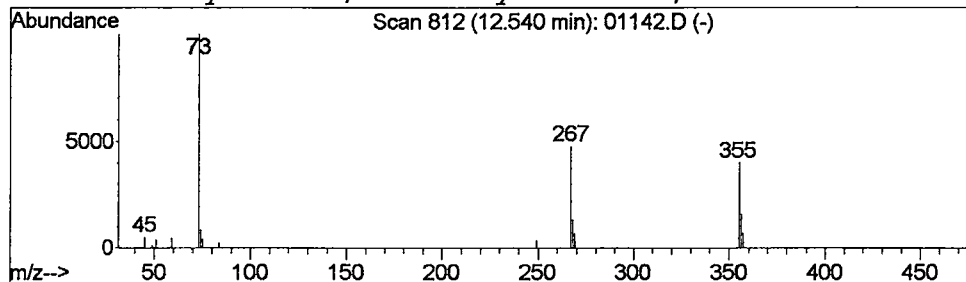
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 Inst : Instrumen
 Multiplr: 1.00

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

 Peak Number 9 Cyclopentasiloxane, decamet... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.54	0.10 ug/L	62986	Naphthalene-d8	13.20

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Cyclopentasiloxane, decamethyl-	370	C10H30O5Si5	000541-02-6	83
2		Cyclopentasiloxane, decamethyl-	370	C10H30O5Si5	000541-02-6	64
3		3,4-DIHYDROXYMANDELIC ACID-TETRATMS	472	C20H40O5Si4	000000-00-0	37
4		Phenethylamine, N-methyl-.beta.,...	399	C18H37NO3Si3	010538-85-9	37



Data File : C:\MSDCHEM\1\5973N\02FEB25\01142.D
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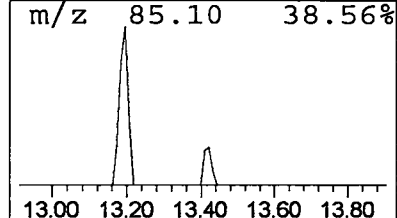
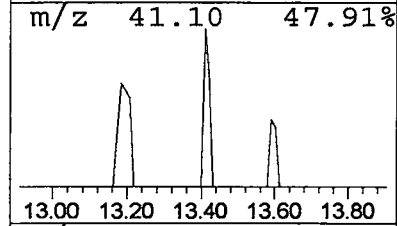
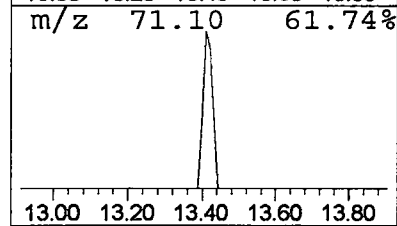
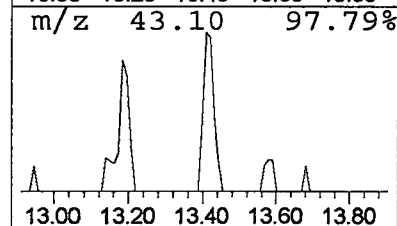
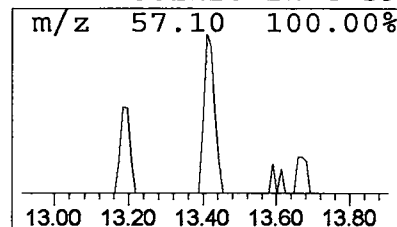
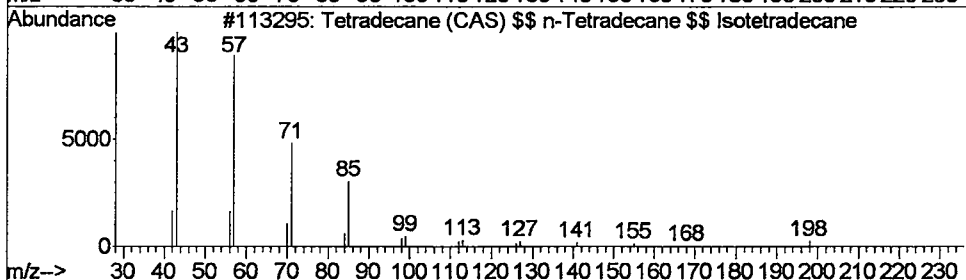
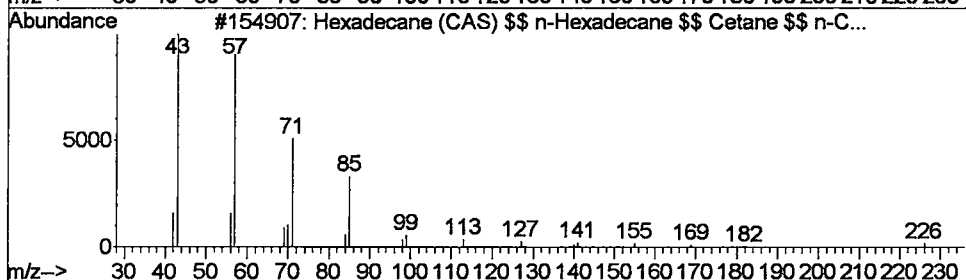
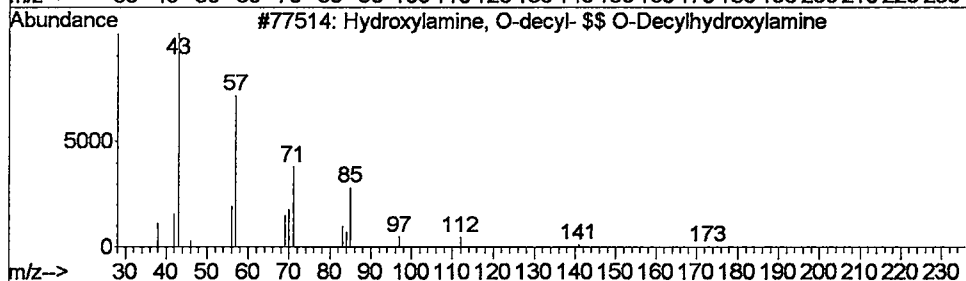
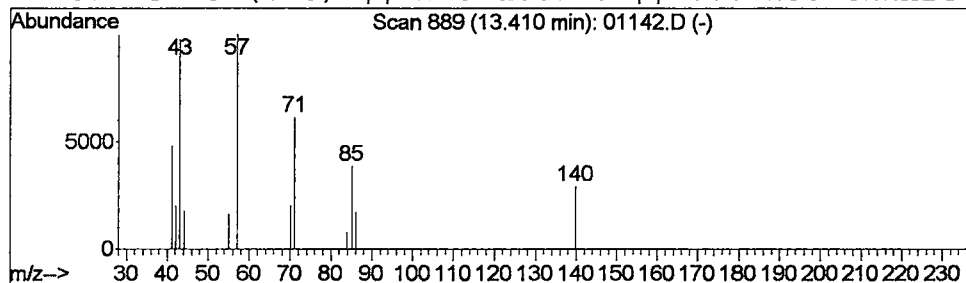
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Quant Method : C:\MSDCHEM\1...\HP022102.M (RTE Integrator)
 Title : 508 scan method
 Library : C:\DATABASE\WILEY7N.L

 Peak Number 10 Hydroxylamine, O-decyl- \$\$... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.41	0.08 ug/L	50520	Naphthalene-d8	13.20

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Hydroxylamine, O-decyl- \$\$ O-Dec...	173	C10H23NO	029812-79-1	59
2			Hexadecane (CAS) \$\$ n-Hexadecane...	226	C16H34	000544-76-3	59
3			Tetradecane (CAS) \$\$ n-Tetradeca...	198	C14H30	000629-59-4	59
4			Undecane (CAS) \$\$ n-Undecane \$\$...	156	C11H24	001120-21-4	59



Data File : C:\MSDCHEM\1\5973N\02FEB25\01142.D
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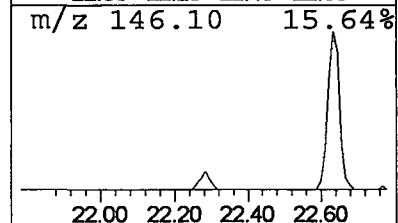
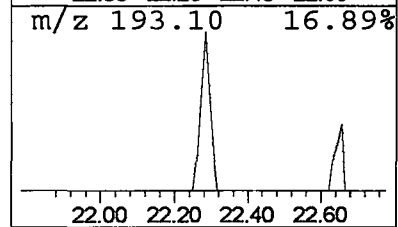
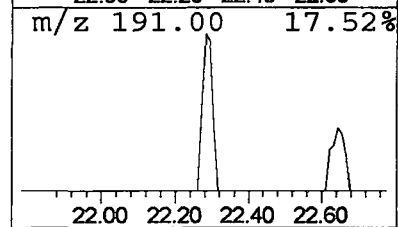
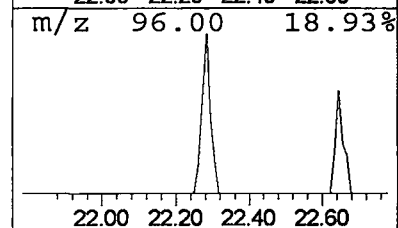
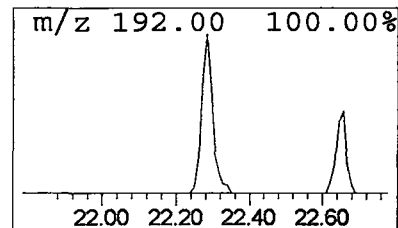
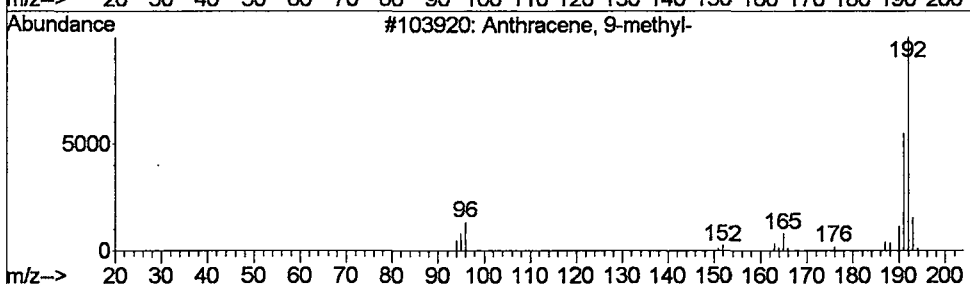
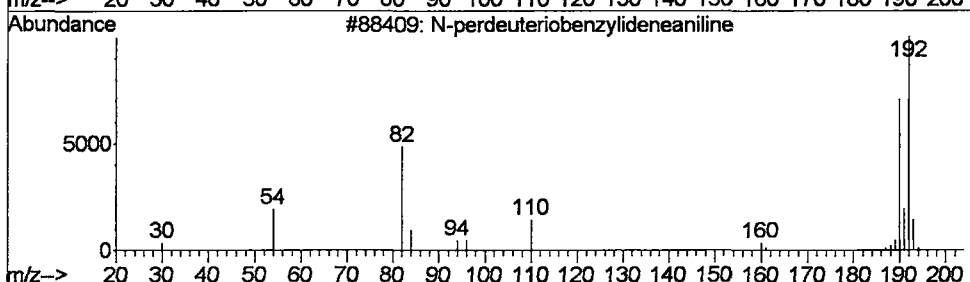
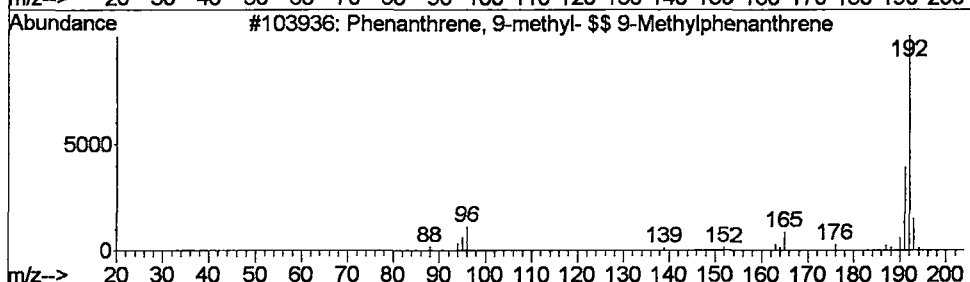
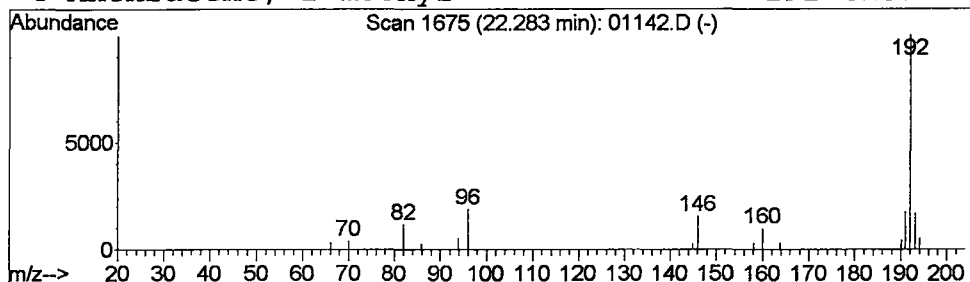
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 Inst : Instrumen
 Multiplr: 1.00

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

 Peak Number 11 Phenanthrene, 9-methyl- \$\$... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.28	0.13 ug/L	89300	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-perdeuteriobenzylideneaniline	192	C13D11N	000538-51-2	58
3		Anthracene, 9-methyl-	192	C15H12	000779-02-2	45
4		Anthracene, 2-methyl-	192	C15H12	000613-12-7	45



Data File : C:\MSDCHEM\1\5973N\02FEB25\01142.D
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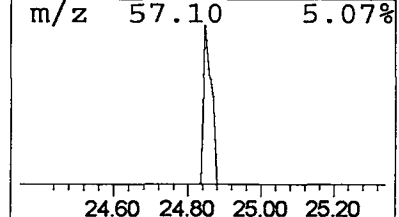
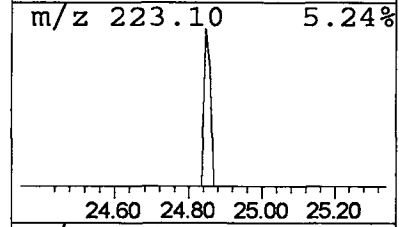
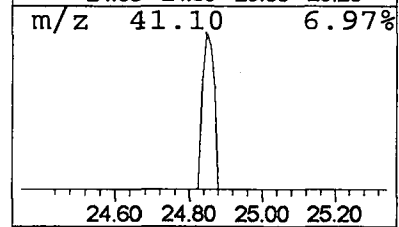
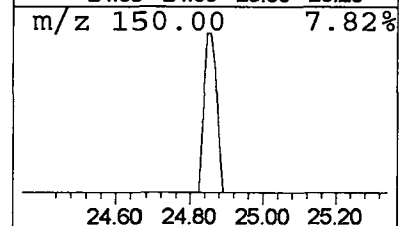
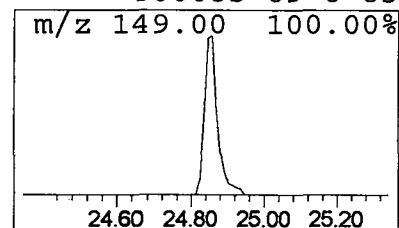
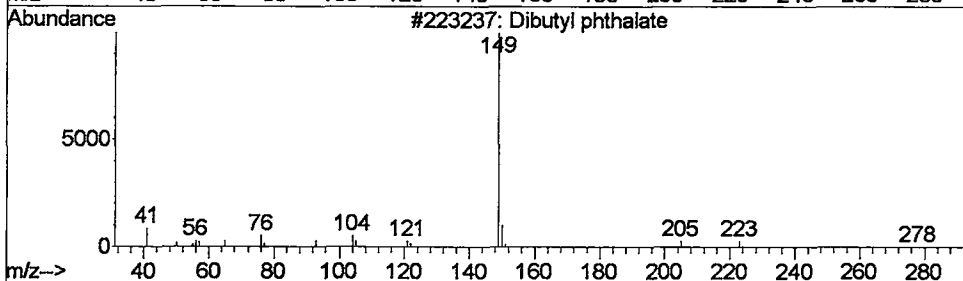
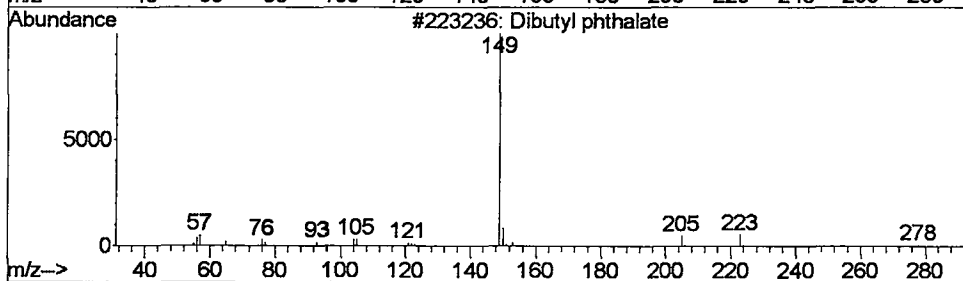
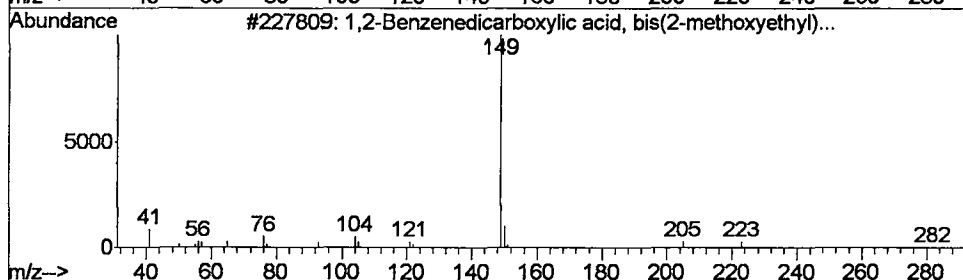
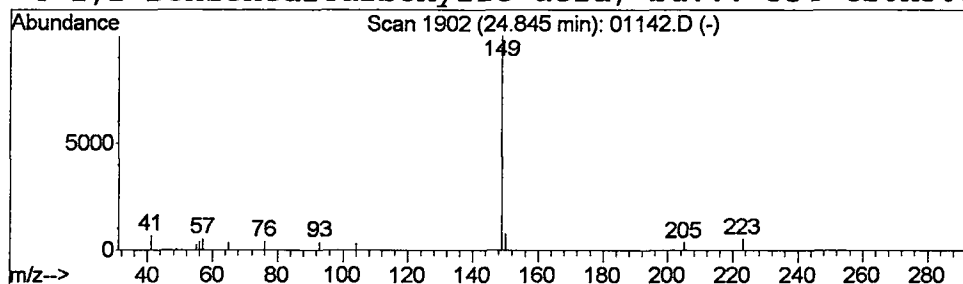
Vial: 8
 Operator:
 Inst : Instrumen
 Multiplr: 1.00

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

 Peak Number 12 1,2-Benzenedicarboxylic aci... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.85	0.12 ug/L	79361	Phenanthrene-d10	22.63

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		1,2-Benzenedicarboxylic acid, bi...	282	C14H18O6	000117-82-8	83
2		Dibutyl phthalate	278	C16H22O4	000084-74-2	83
3		Dibutyl phthalate	278	C16H22O4	000084-74-2	83
4		1,2-Benzenedicarboxylic acid, bu...	334	C20H30O4	000085-69-8	83



Data File : C:\MSDCHEM\1\5973N\02FEB25\01142.D
 Acq On : 25 Feb 2002 5:01 pm
 Sample : 508 scan
 Misc : February 25, 2002
 MS Integration Params: LSCINT.P

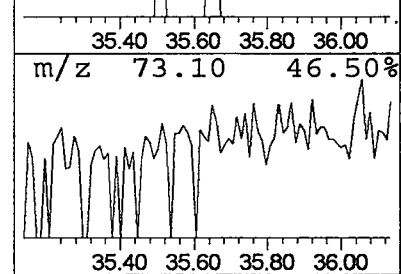
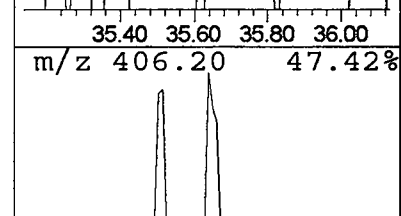
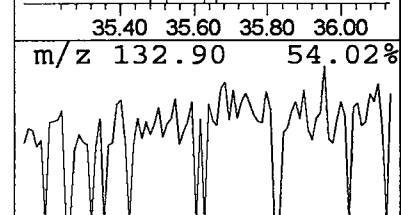
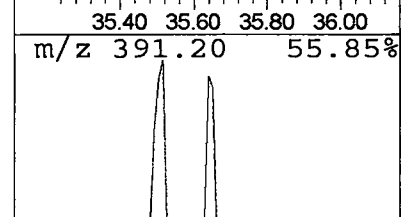
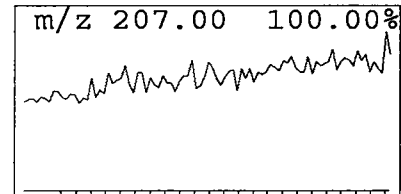
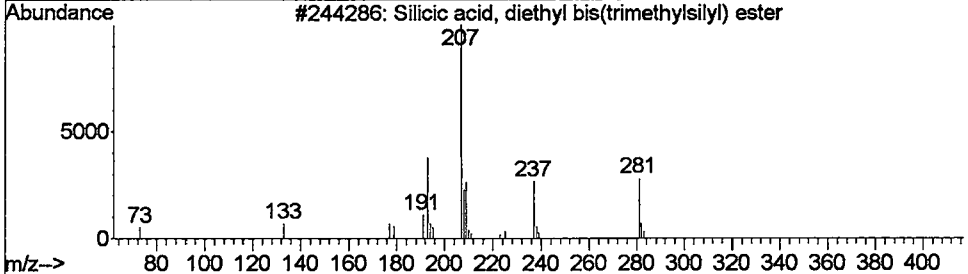
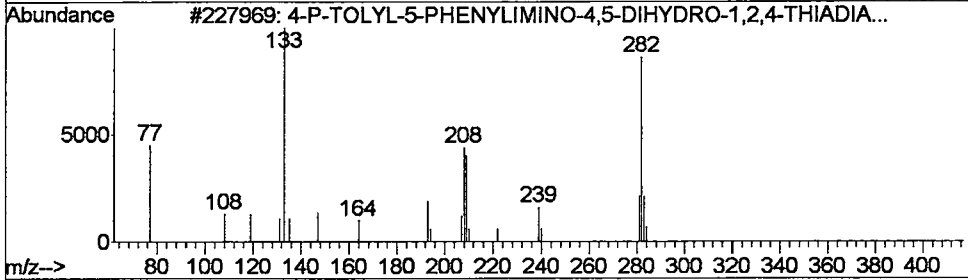
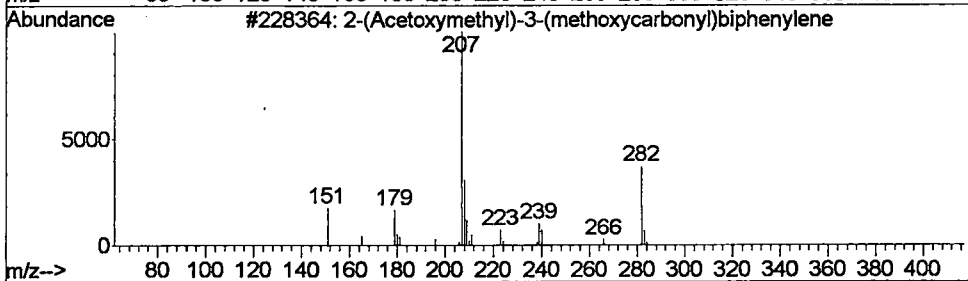
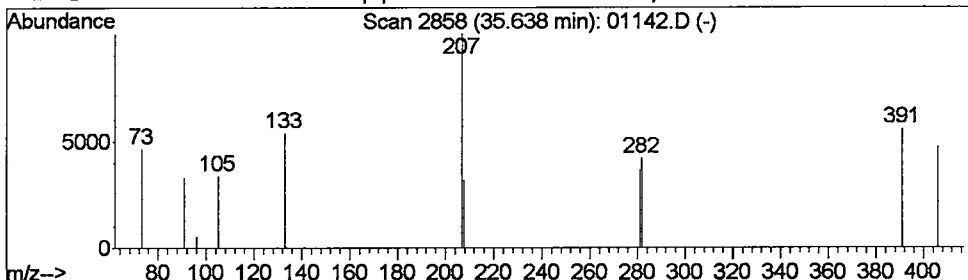
Vial: 8
 Operator:
 Inst : Instrumen
 Multiplr: 1.00

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

 Peak Number 13 2-(Acetoxymethyl)-3-(methox... Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
35.64	0.05 ug/L	24017	Perylene-d12	34.41

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-(Acetoxymethyl)-3-(methoxycarb...	282	C17H14O4	093103-70-9	32
2			4-P-TOLYL-5-PHENYLIMINO-4,5-DIHY...	282	C15H14N4S	054568-21-7	16
3			Silicic acid, diethyl bis(trimet...	296	C10H28O4Si3	003555-45-1	9
4			Gibberellin A3 \$\$ Gibb-3-ene-1,1...	346	C19H22O6	000077-06-5	9



Operator ID: Date Acquired: 25 Feb 2002 5:01 pm
 Data File: C:\MSDCHEM\1\5973N\02FEB25\01142.D
 Name: 508 scan
 Misc: February 25, 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
Dichloroacetoneitr...	3.50	0.2	ug/L	73427	1	9.72	8534010	20.0
Butanoic acid, me...	3.62	0.1	ug/L	60193	1	9.72	8534010	20.0
Propane, 1,1-dime...	3.94	0.1	ug/L	32748	1	9.72	8534010	20.0
Benzene, methyl- ...	4.39	0.1	ug/L	48469	1	9.72	8534010	20.0
3-(CYCLOHEX-3'-EN...	5.94	0.3	ug/L	137872	1	9.72	8534010	20.0
2-Propanone, 1,1,...	6.00	0.3	ug/L	138635	1	9.72	8534010	20.0
3-Butenoic acid \$...	9.90	0.1	ug/L	28692	1	9.72	8534010	20.0
4-Cyanocyclohexen...	9.96	0.1	ug/L	26666	1	9.72	8534010	20.0
Cyclopentasiloxan...	12.54	0.1	ug/L	62986	2	13.20	12350700	20.0
Hydroxylamine, O-...	13.41	0.1	ug/L	50520	2	13.20	12350700	20.0
Phenanthrene, 9-m...	22.28	0.1	ug/L	89300	4	22.63	13338100	20.0
1,2-Benzenedicarb...	24.85	0.1	ug/L	79361	4	22.63	13338100	20.0
2-(Acetoxymethyl)...	35.64	0.0	ug/L	24017	6	34.41	9977560	20.0