

User: Vinci, David L
Westchester Dept. of Labs & Research
Date: 03/05/2002 Time: 12:15:26

Sample: AE02717
Analysis: \$GCMS GCMS Scan For Unknowns
Units: ug
Started: 3/5/02 12:14 Ended: 3/5/02 12:14 Analyst: DLV
Page: 1

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

Comments for Sample AE02717 - Analysis \$GCMS

Westchester Dept. of Labs & Research

User: Vinci, David L

Date: 03-05-2002 Time: 12:15:50

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

Quantitation Report (Not Reviewed)

Data File : C:\MSDCHEM\1\5973N\02FEB28\02717.D Vial: 2
 Acq On : 28 Feb 2002 3:30 pm Operator:
 Sample : 508 scan Inst : Instrumen
 Misc : February 28, 2002 Multiplr: 1.00
 MS Integration Params: SPLIT.P
 Quant Time: Mar 01 09:02:55 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	1313527	20.00	ug/L	0.00
10) Naphthalene-d8	13.20	136	5584922	20.00	ug/L	0.00
17) Acenaphthene-d10	18.33	164	2758535	20.00	ug/L	0.00
29) Phenanthrene-d10	22.63	188	4894654	20.00	ug/L	-0.01
38) Chrysene-d12	30.49	240	3805823	20.00	ug/L	-0.03
44) Perylene-d12	34.41	264	2492773	20.00	ug/L	-0.04

System Monitoring Compounds

37) 2,4-DDD	27.72	235	284096	4.35	ug/L	0.00
Spiked Amount	5.000		Recovery	=	87.00%	

Target Compounds

	R.T.	QIon	Response	Conc	Units	Qvalue
20) Dimethylphthalate	18.34	163	448722	2.68	ug/L #	58
21) 2,6-Dinitrotoluene	18.33	165	347769	9.00	ug/L #	27

Quantitation Report

(Not Reviewed)

Data File : C:\MSDCHEM\1\5973N\02FEB28\02717.D

Vial: 2

Acq On : 28 Feb 2002 3:30 pm

Operator:

Sample : 508 scan

Inst : Instrumen

Misc : February 28, 2002

Multiplr: 1.00

MS Integration Params: SPLIT.P

Quant Time: Mar 1 9:02 2002

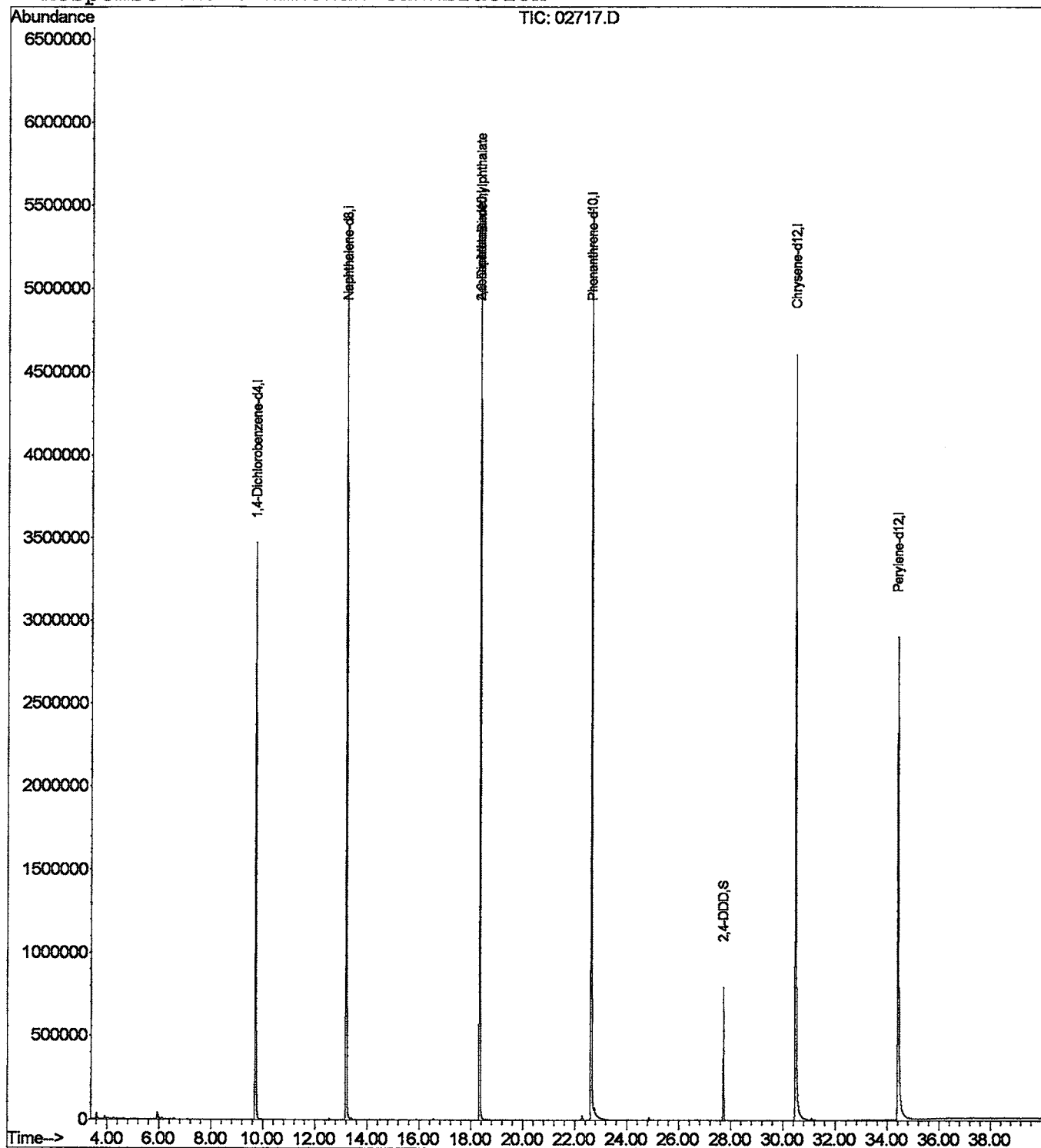
Quant Results File: HP022102.R

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



02717.D HP022102.M

Fri Mar 01 09:02:57 2002

Page 2

LSC Area Percent Report

Data File : C:\MSDCHEM\1\5973N\02FEB28\02717.D
 Acq On : 28 Feb 2002 3:30 pm
 Sample : 508 scan
 Misc : February 28, 2002
 MS Integration Params: LSCINT.P

Vial: 2
 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

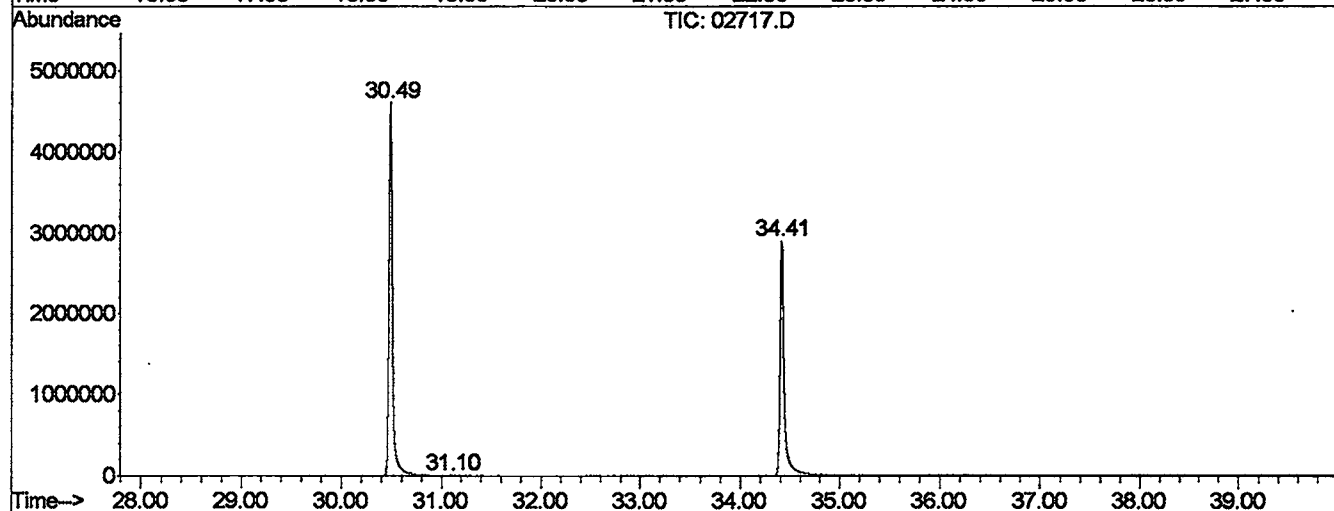
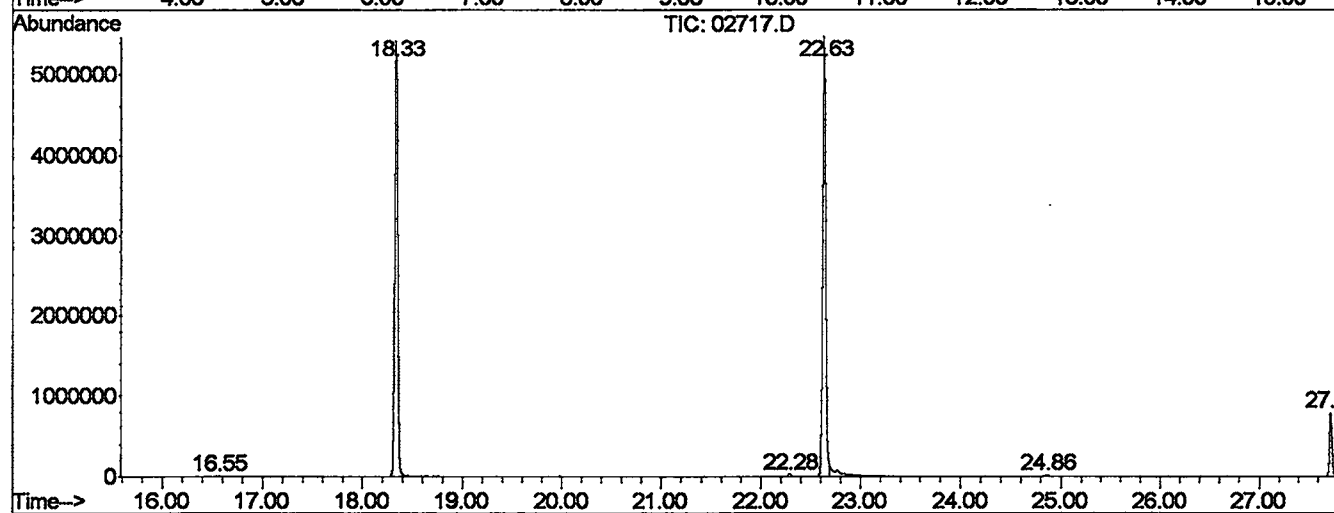
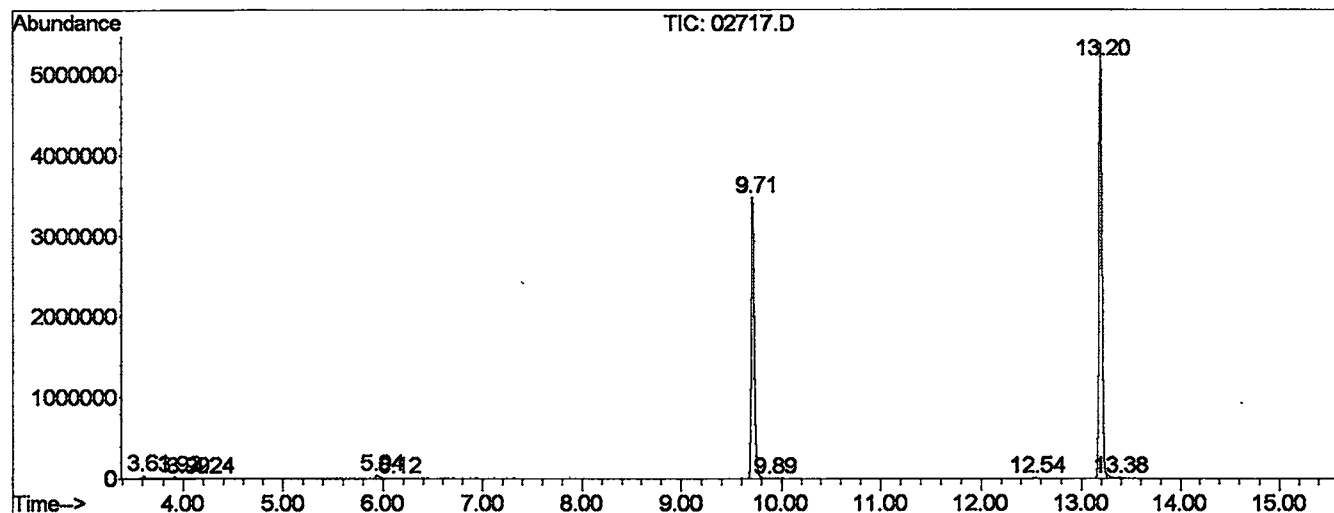
Signal : TIC

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.611	19	21	27	rBV	35421	56700	0.47%	0.091%
2	3.916	45	48	53	rBV	19708	27400	0.23%	0.044%
3	3.995	53	55	61	rBV4	4811	17400	0.15%	0.028%
4	4.243	75	77	85	rVV5	6332	17474	0.15%	0.028%
5	5.936	223	227	239	rBV	42412	126559	1.06%	0.203%
6	6.117	239	243	259	rVB4	7073	29490	0.25%	0.047%
7	9.707	557	561	575	rVV	3474180	7318692	61.25%	11.732%
8	9.888	575	577	595	rVB4	4093	30498	0.26%	0.049%
9	12.540	807	812	815	rBV	8357	17488	0.15%	0.028%
10	13.195	865	870	875	rBV	5381935	10476656	87.67%	16.794%
11	13.376	885	886	893	rVB	11694	20438	0.17%	0.033%
12	16.548	1163	1167	1171	rBV	9819	18866	0.16%	0.030%
13	18.332	1319	1325	1331	rBV	5425195	11175695	93.52%	17.914%
14	22.283	1671	1675	1681	rVV	29875	62034	0.52%	0.099%
15	22.633	1701	1706	1711	rBV2	5475567	11949554	100.00%	19.155%
16	24.857	1899	1903	1915	rVV	16816	38545	0.32%	0.062%
17	27.724	2153	2157	2163	rBV	792984	1374894	11.51%	2.204%
18	30.490	2397	2402	2439	rBV2	4604381	11147243	93.29%	17.869%
19	31.099	2451	2456	2461	rVB2	12079	23756	0.20%	0.038%
20	34.407	2743	2749	2781	rBV	2895021	8455471	70.76%	13.554%

Sum of corrected areas: 62384853

LSC Report - Integrated Chromatogram

File : C:\MSDCHEM\1\5973N\02FEB28\02717.D
 Operator :
 Acquired : 28 Feb 2002 3:30 pm using AcqMethod HP020702
 Instrument : Instrumen
 Sample Name: 508 scan
 Misc Info : February 28, 2002
 Vial Number: 2
 Quant File :HP022102.RES (RTE Integrator)



Library Search Compound Report

Data File : C:\MSDCHEM\1\5973N\02FEB28\02717.D
 Acq On : 28 Feb 2002 3:30 pm
 Sample : 508 scan
 Misc : February 28, 2002
 MS Integration Params: LSCINT.P

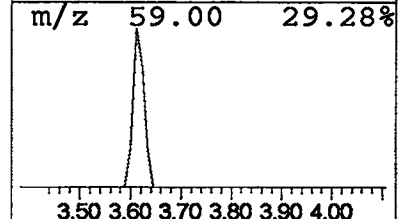
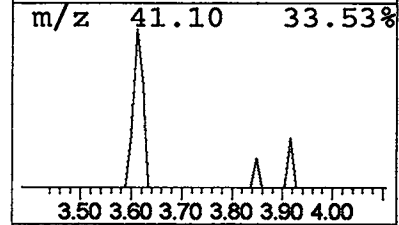
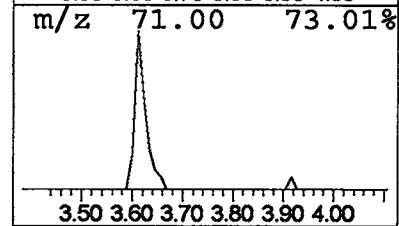
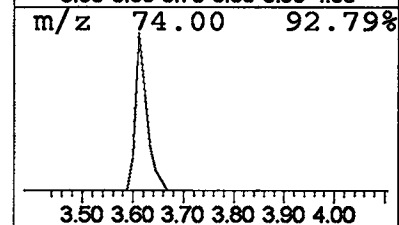
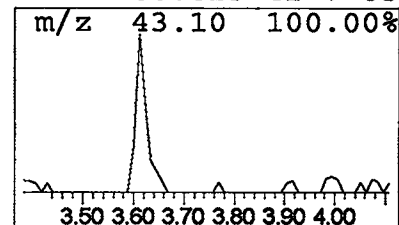
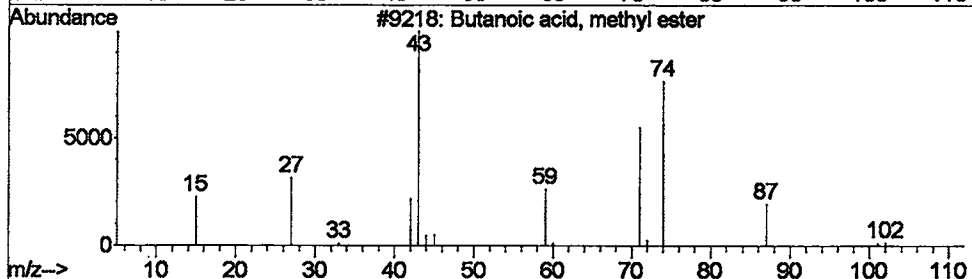
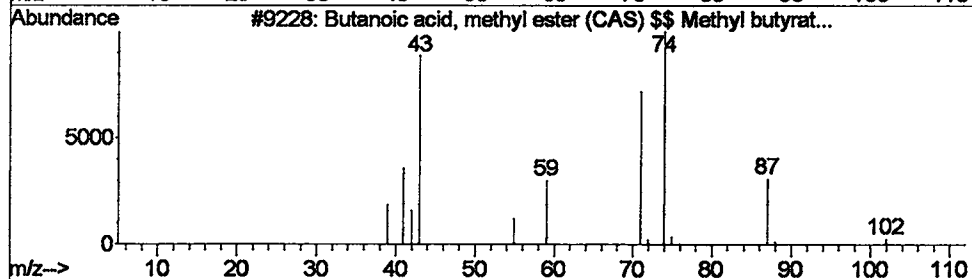
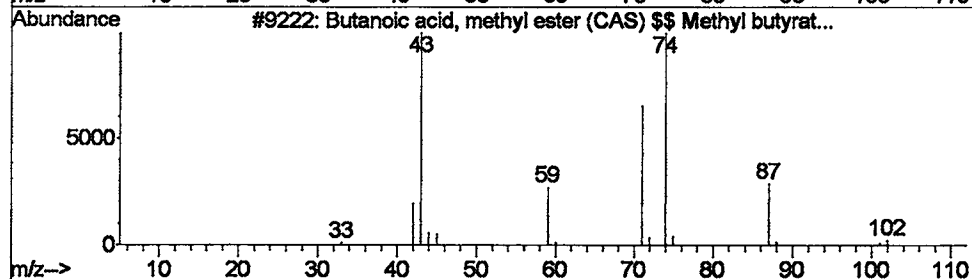
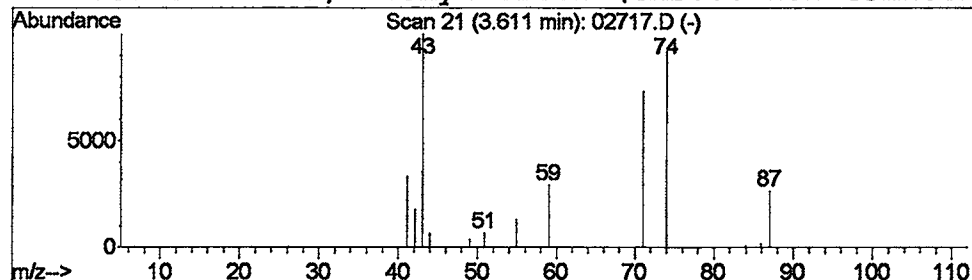
Vial: 2
 Operator:
 Inst : Instrumen
 Multiplr: 1.00

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

 Peak Number 1 Butanoic acid, methyl ester... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.61	0.15 ug/L	56700	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\02FEB28\02717.D
 Acq On : 28 Feb 2002 3:30 pm
 Sample : 508 scan
 Misc : February 28, 2002
 MS Integration Params: LSCINT.P

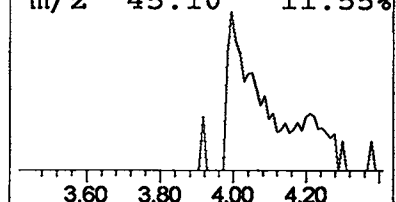
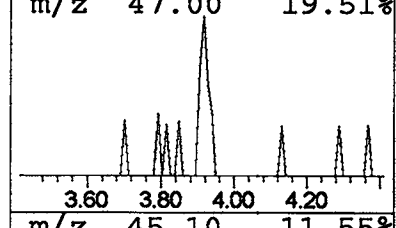
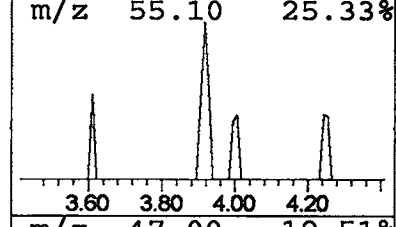
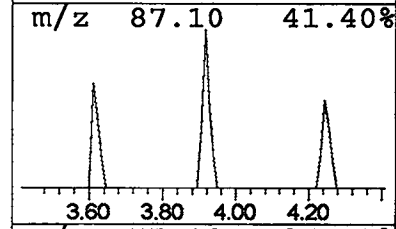
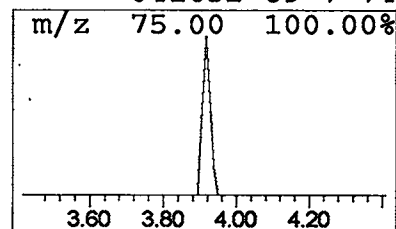
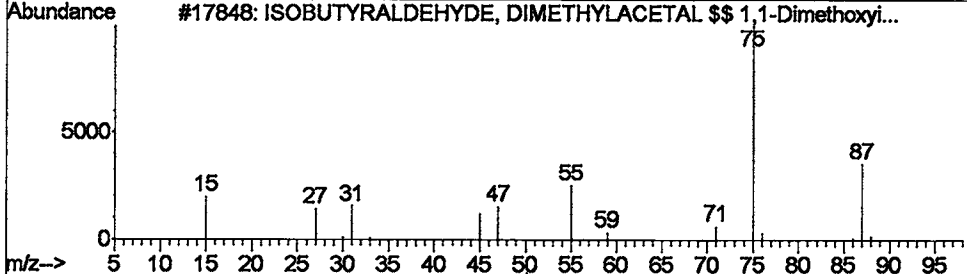
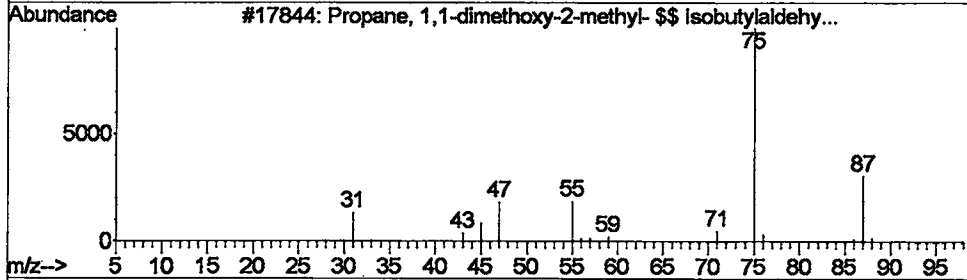
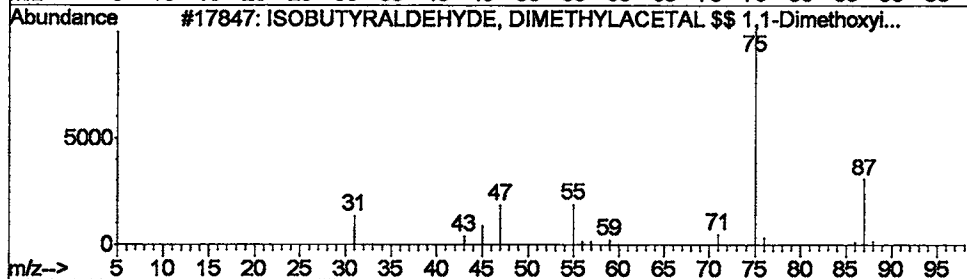
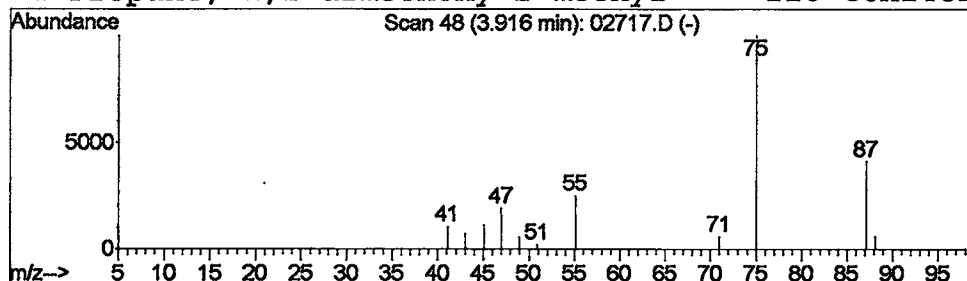
Vial: 2
 Operator:
 Inst : Instrumen
 Multiplr: 1.00

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

 Peak Number 2 ISOBUTYRALDEHYDE, DIMETHYLA... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.92	0.07 ug/L	27400	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			Propane, 1,1-dimethoxy-2-methyl-	118	C6H14O2	041632-89-7	74



Library Search Compound Report

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Acq On : 28 Feb 2002 3:30 pm
Sample : 508 scan
Misc : February 28, 2002
MS Integration Params: LSCINT.P

Vial: 2
Operator:
Inst : Instrumen
Multiplr: 1.00

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

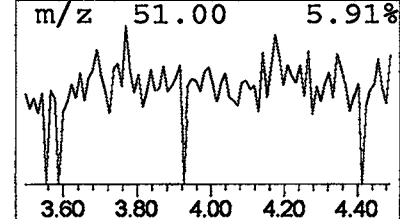
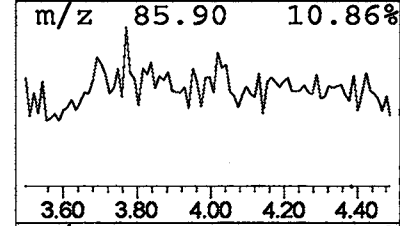
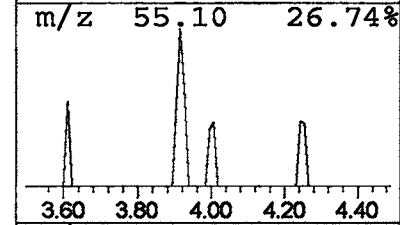
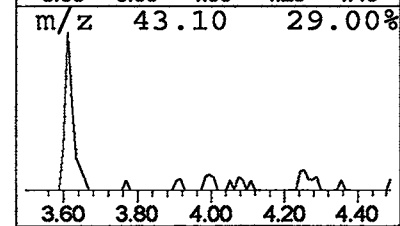
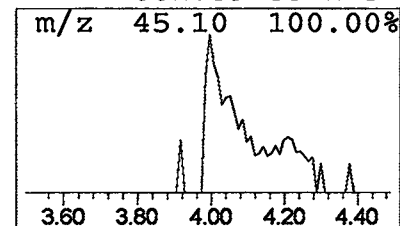
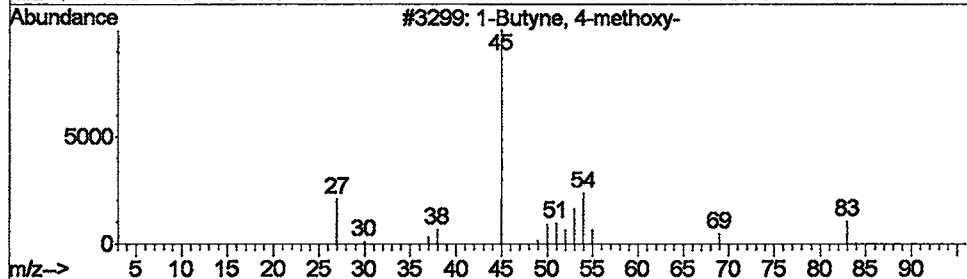
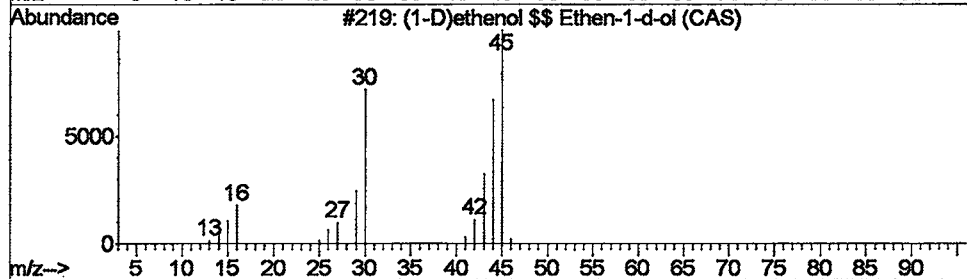
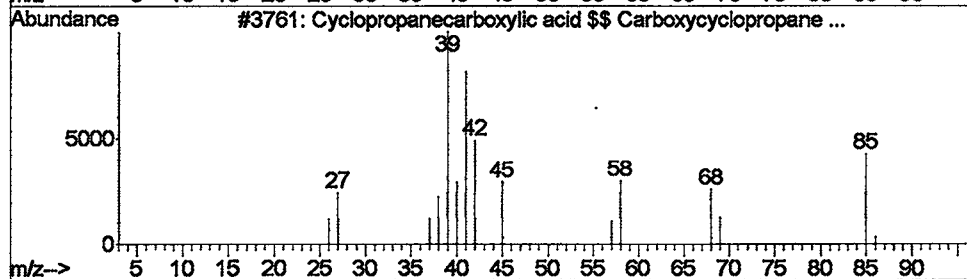
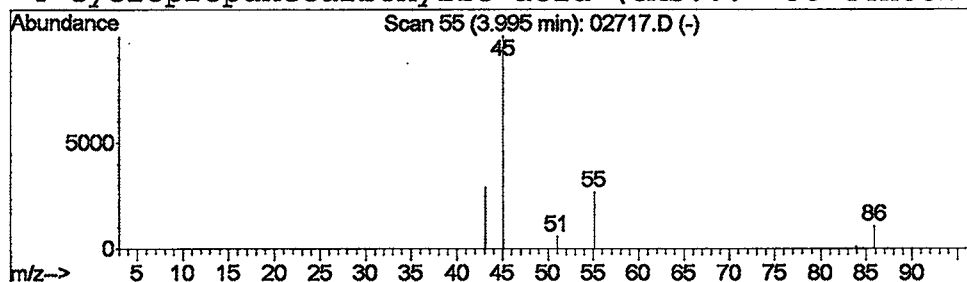
Title : 508 scan method

Library : C:\DATABASE\WILEY7N.L

Peak Number 3 Cyclopropanecarboxylic acid... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.99	0.05 ug/L	17400	1,4-Dichlorobenzene-d4	9.72

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclopropanecarboxylic acid \$\$ C...	86	C4H6O2	001759-53-1	3
2			(1-D)ethenol \$\$ Ethen-1-d-ol (CAS)	45	C2H3DO	090134-16-0	3
3			1-Butyne, 4-methoxy-	84	C5H8O	036678-08-7	3
4			Cyclopropanecarboxylic acid (CAS...	86	C4H6O2	001759-53-1	3



Library Search Compound Report

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 MS Integration Params: LSCINT.P

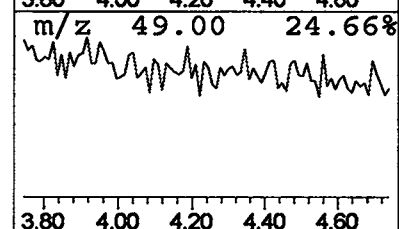
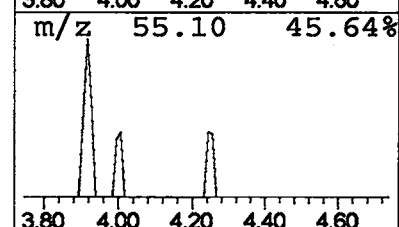
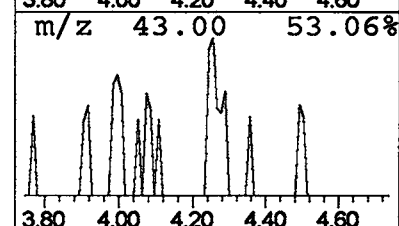
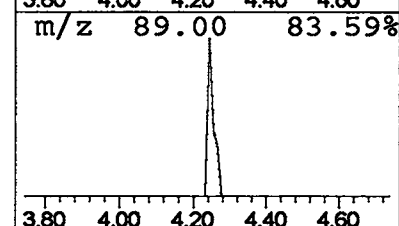
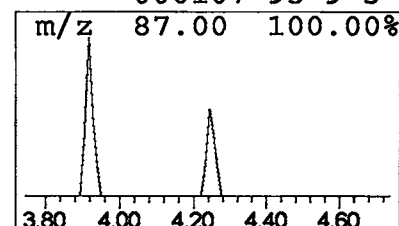
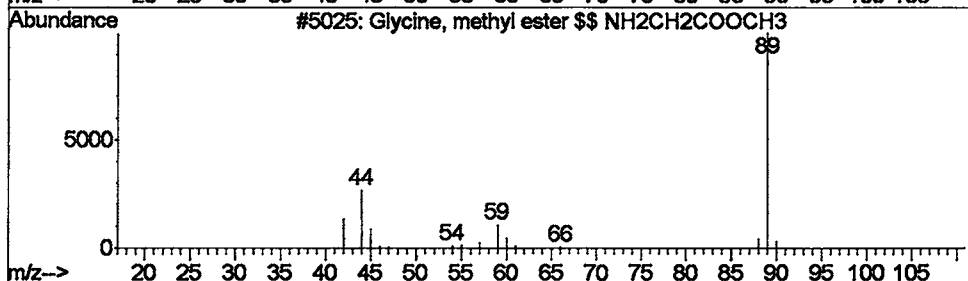
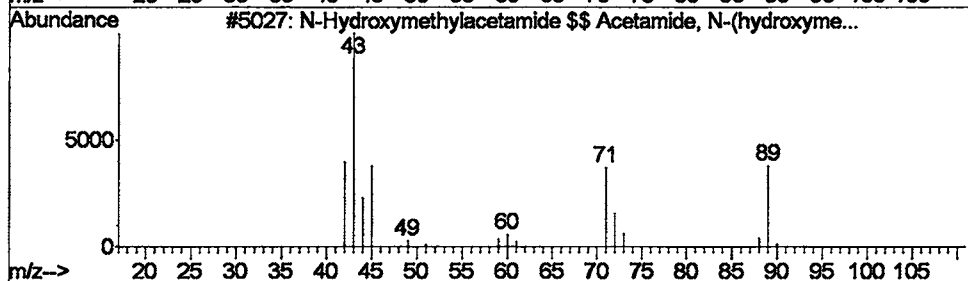
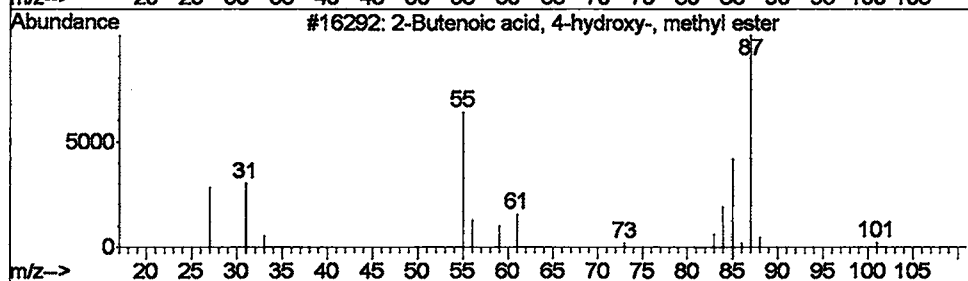
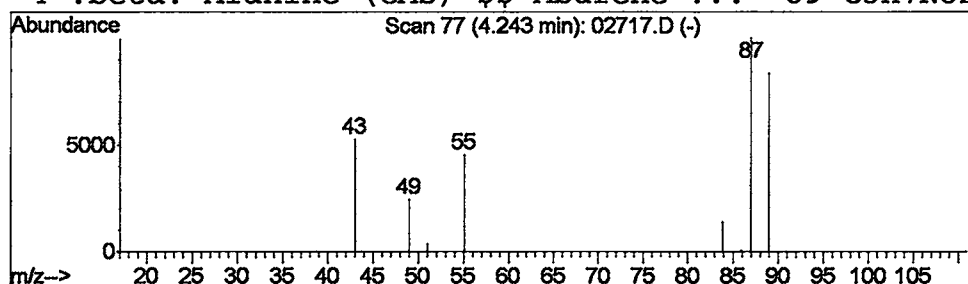
Vial: 2
 Operator:
 Inst : Instrumen
 Multiplr: 1.00

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

 Peak Number 4 2-Butenoic acid, 4-hydroxy-... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.24	0.05 ug/L	17474	1,4-Dichlorobenzene-d4	9.72

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Butenoic acid, 4-hydroxy-, met...	116	C5H8O3	004508-99-0	4
2			N-Hydroxymethylacetamide \$\$ Acet...	89	C3H7NO2	000625-51-4	3
3			Glycine, methyl ester \$\$ NH2CH2C...	89	C3H7NO2	000616-34-2	3
4			.beta.-Alanine (CAS) \$\$ Abufene ...	89	C3H7NO2	000107-95-9	3



Library Search Compound Report

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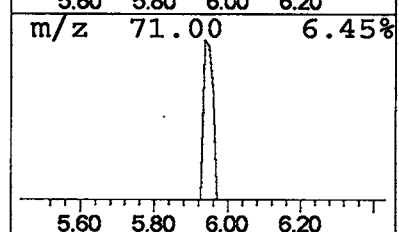
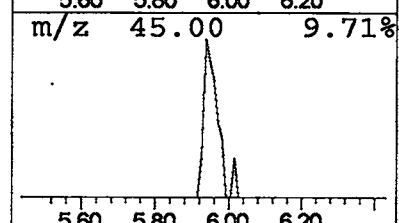
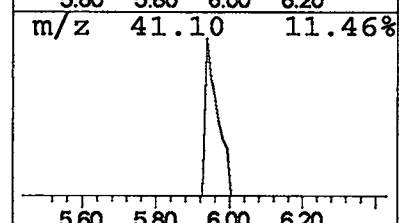
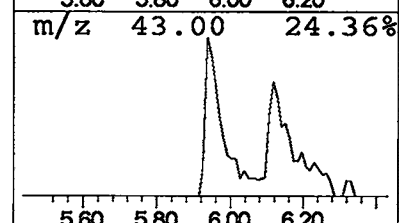
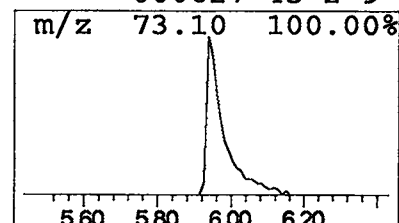
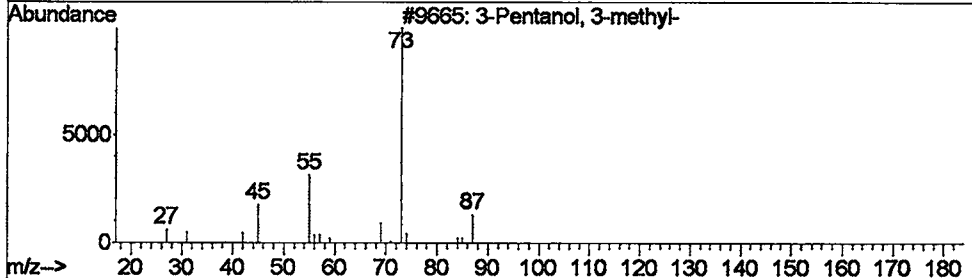
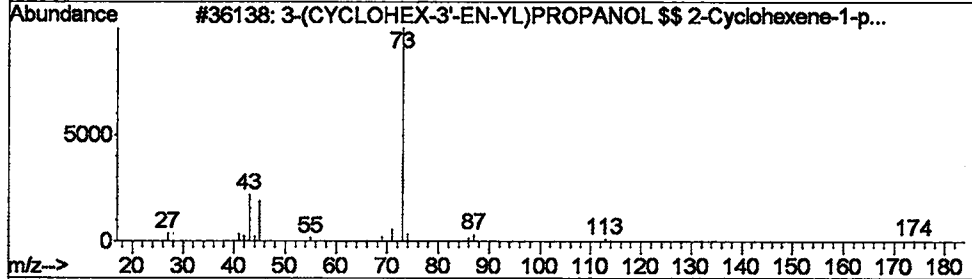
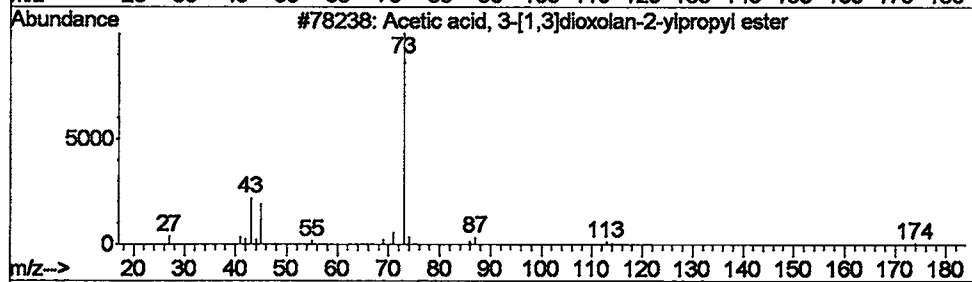
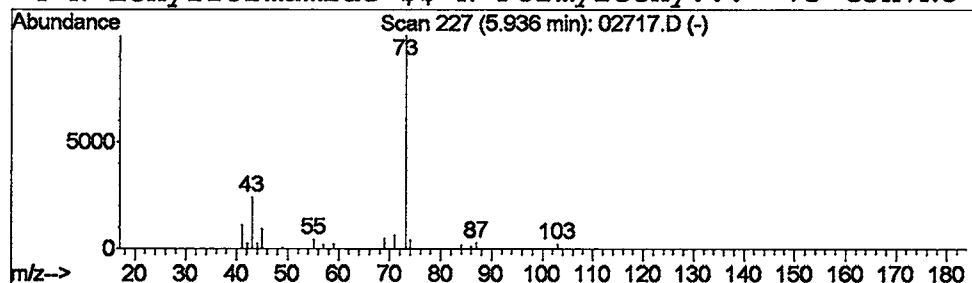
Vial: 2
 Operator:
 Inst : Instrumen
 Multiplr: 1.00

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

 Peak Number 5 Acetic acid, 3-[1,3]dioxola... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.94	0.35 ug/L	126559	1,4-Dichlorobenzene-d4	9.72

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Acetic acid, 3-[1,3]dioxolan-2-y...	174	C8H14O4	000000-00-0	33	
2	3-(CYCLOHEX-3'-EN-YL)PROPANOL \$\$...	174	C8H14O4	015745-87-6	33	
3	3-Pentanol, 3-methyl-	102	C6H14O	000077-74-7	28	
4	N-Ethylformamide \$\$ N-Formylethy...	73	C3H7NO	000627-45-2	9	



Library Search Compound Report

Data File : C:\MSDCHEM\1\5973N\02FEB28\02717.D
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Sample : 508 scan
Misc : February 28, 2002
MS Integration Params: LSCINT.P

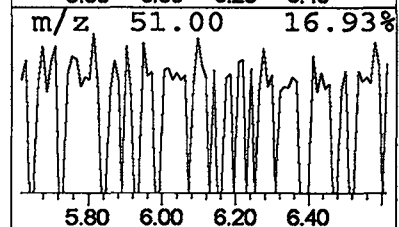
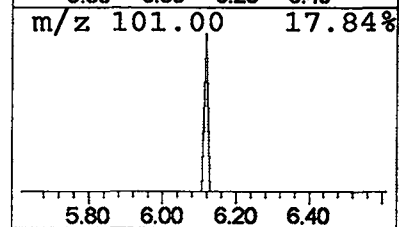
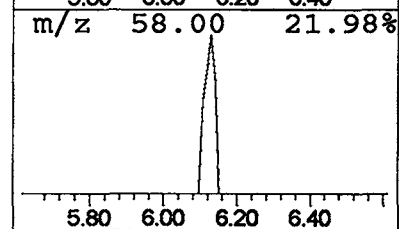
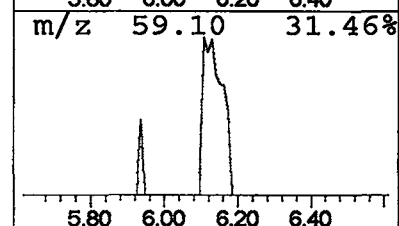
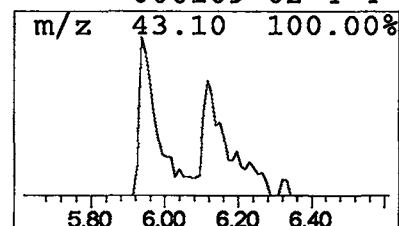
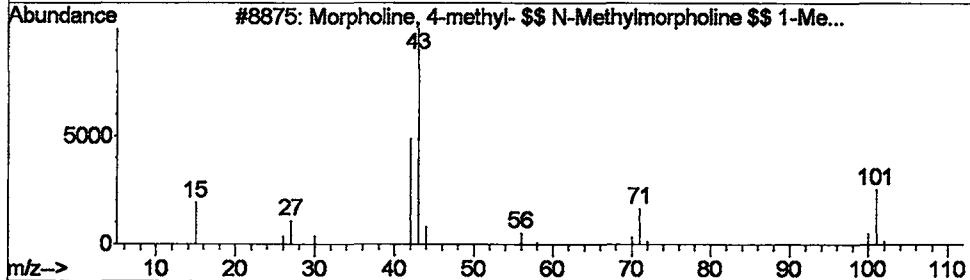
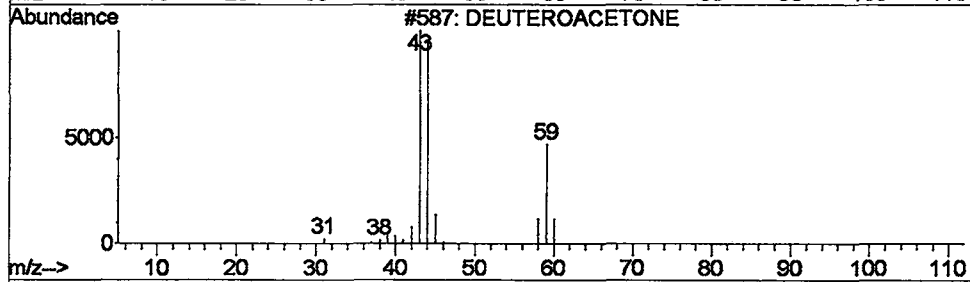
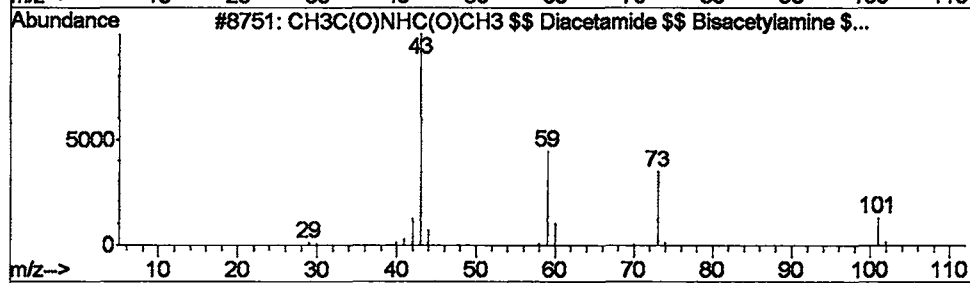
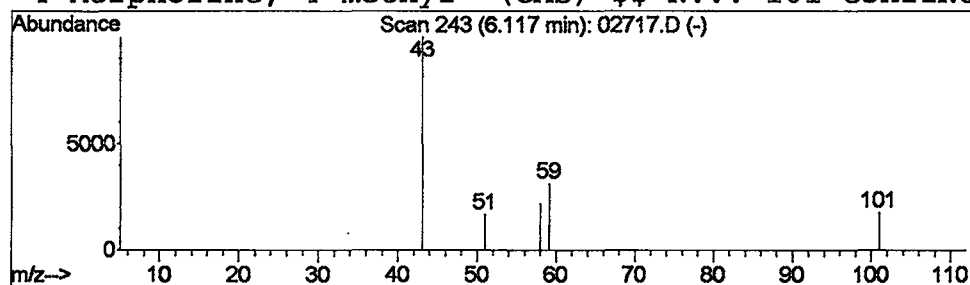
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Inst : Instrumen
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Title : 508 scan method
Library : C:\DATABASE\WILEY7N.L

Peak Number 6 CH3C(O)NHC(O)CH3 \$\$ Diaceta... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.12	0.08 ug/L	29490	1,4-Dichlorobenzene-d4	9.72

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		CH3C(O)NHC(O)CH3 \$\$ Diacetamide ...	101	C4H7NO2	000625-77-4	7
2		DEUTEROACETONE	59	C3H5DO	004468-52-4	4
3		Morpholine, 4-methyl- \$\$ N-Methy...	101	C5H11NO	000109-02-4	4
4		Morpholine, 4-methyl- (CAS) \$\$ N...	101	C5H11NO	000109-02-4	4



Library Search Compound Report

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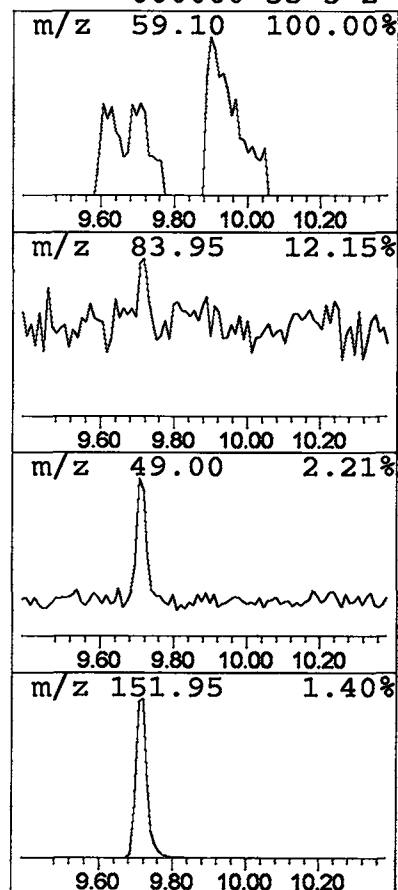
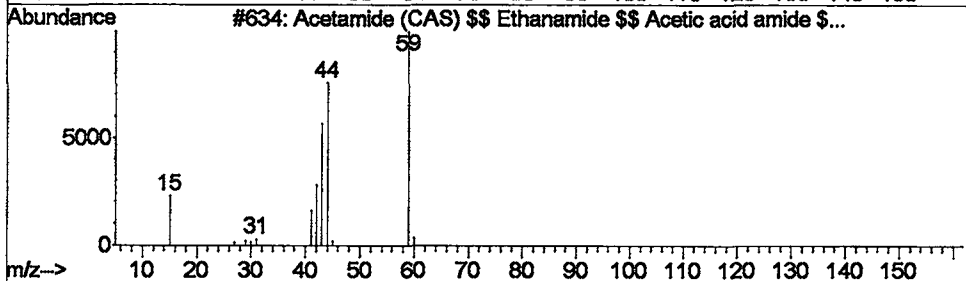
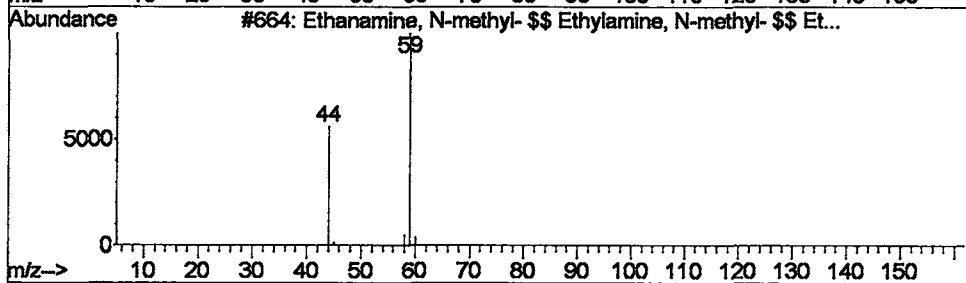
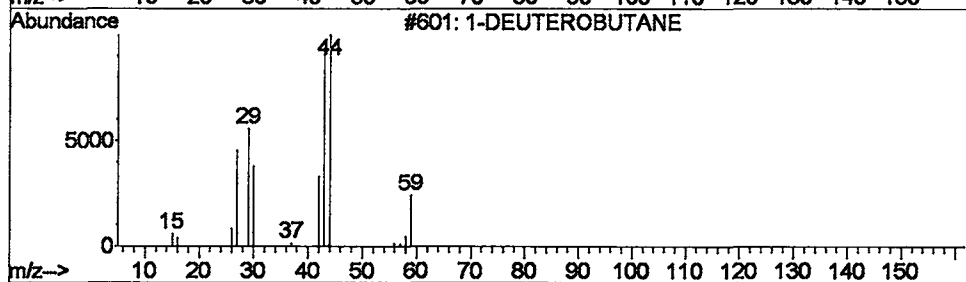
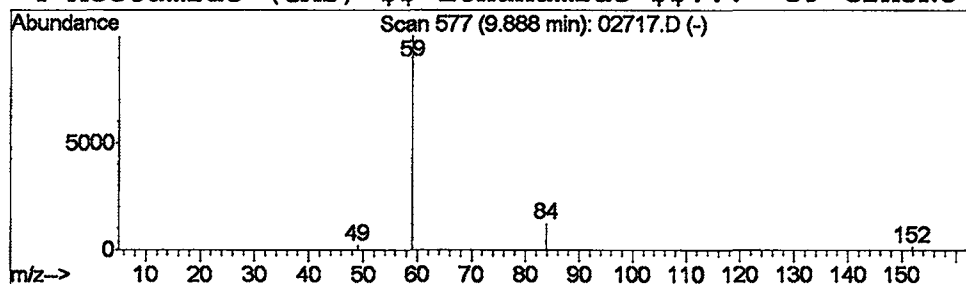
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Quant Method : C:\MSDCHEM\1...\HP022102.M (RTE Integrator)
 Title : 508 scan method
 Library : C:\DATABASE\WILEY7N.L

 Peak Number 7 1-DEUTEROBUTANE Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.89	0.08 ug/L	30498	1,4-Dichlorobenzene-d4	9.72

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-DEUTEROBUTANE	59	C4H9D	053716-53-3	3
2	Ethanamine, N-methyl- \$\$ Ethylam...	59	C3H9N	000624-78-2	2
3	Acetamide (CAS) \$\$ Ethanamide \$\$...	59	C2H5NO	000060-35-5	2
4	Acetamide (CAS) \$\$ Ethanamide \$\$...	59	C2H5NO	000060-35-5	2



Library Search Compound Report

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Acq On : 28 Feb 2002 3:30 pm
Sample : 508 scan
Misc : February 28, 2002
MS Integration Params: LSCINT.P

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Multiplr: 1.00

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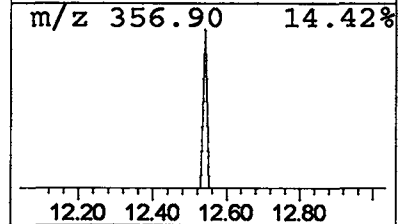
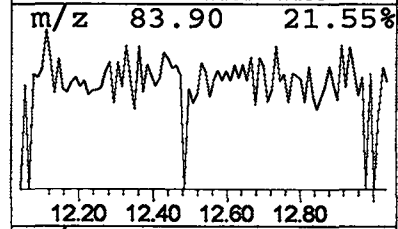
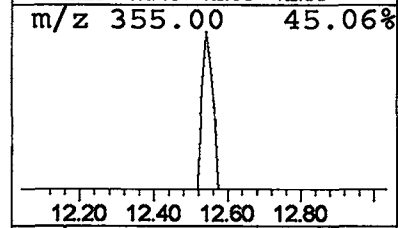
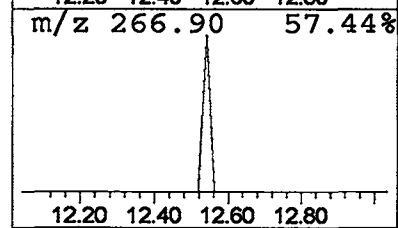
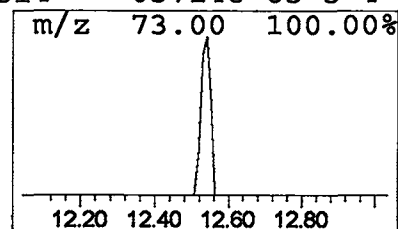
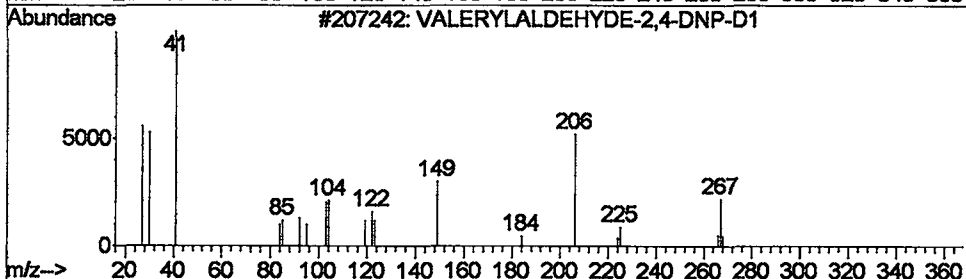
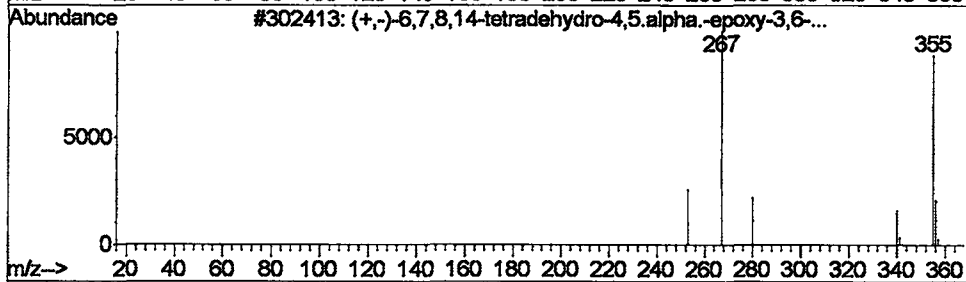
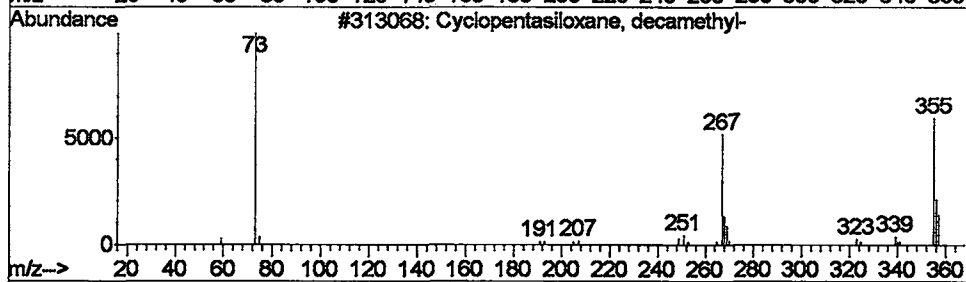
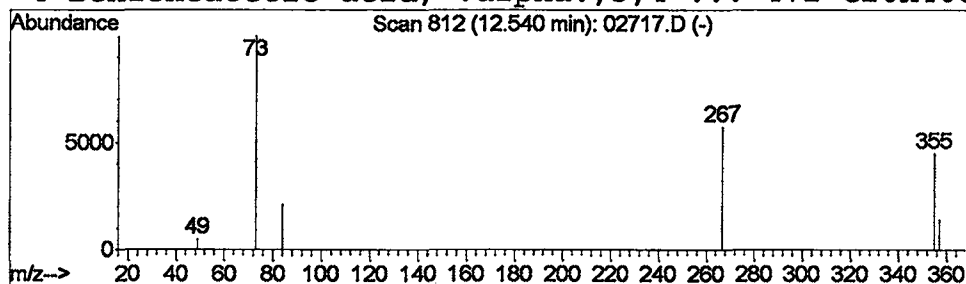
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Library : C:\DATABASE\WILEY7N.L

Peak Number 8 Cyclopentasiloxane, decamet... Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.54	0.03 ug/L	17488	Naphthalene-d8	13.20

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclopentasiloxane, decamethyl-	370	C10H30O5Si5	000541-02-6	39
2			(+,-)-6,7,8,14-tetradecahydro-4,5....	355	C20H21NO5	113948-35-9	7
3			VALERYLALDEHYDE-2,4-DNP-D1	267	C11H13DN4O4	027820-00-4	4
4			Benzeneacetic acid, .alpha.,3,4-...	472	C20H40O5Si4	037148-65-5	4



Library Search Compound Report

Data File : C:\MSDCHEM\1\5973N\02FEB28\02717.D
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Sample : 508 scan
Misc : February 28, 2002
MS Integration Params: LSCINT.P

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Multiplr: 1.00

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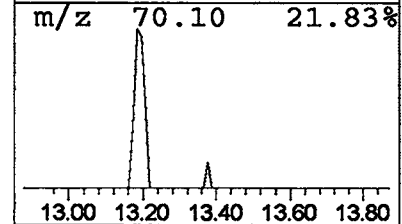
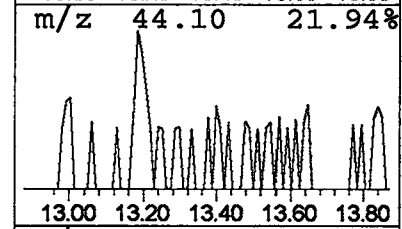
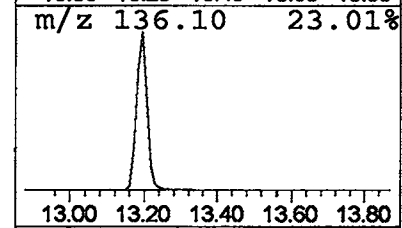
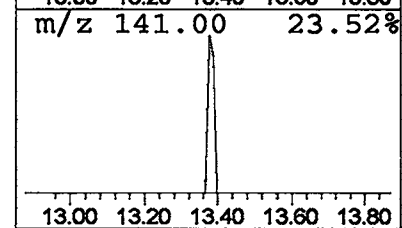
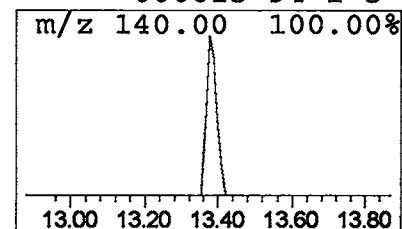
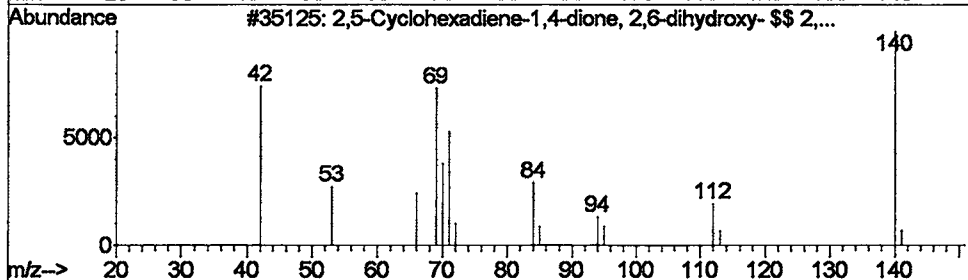
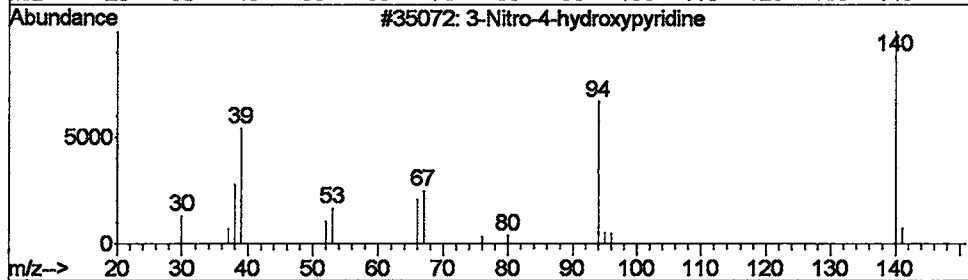
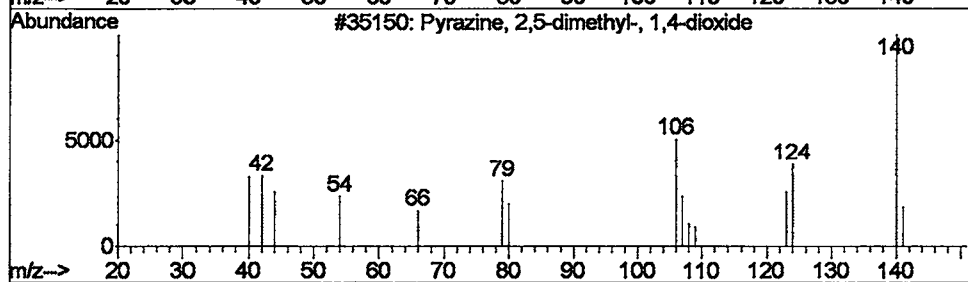
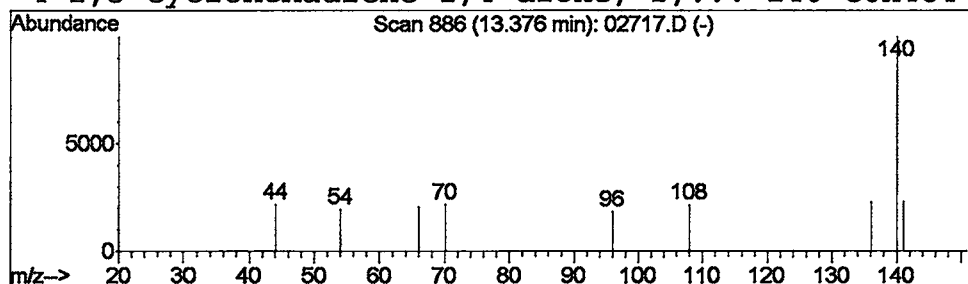
Title : 508 scan method

Library : C:\DATABASE\WILEY7N.L

Peak Number 9 Pyrazine, 2,5-dimethyl-, 1,... Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.38	0.04 ug/L	20438	Naphthalene-d8	13.20

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Pyrazine, 2,5-dimethyl-, 1,4-dio...	140	C6H8N2O2	006890-38-6	45
2			3-Nitro-4-hydroxypyridine	140	C5H4N2O3	000000-00-0	5
3			2,5-Cyclohexadiene-1,4-dione, 2,...	140	C6H4O4	035069-70-6	5
4			2,5-Cyclohexadiene-1,4-dione, 2,...	140	C6H4O4	000615-94-1	5



Library Search Compound Report

Data File : C:\MSDCHEM\1\5973N\02FEB28\02717.D
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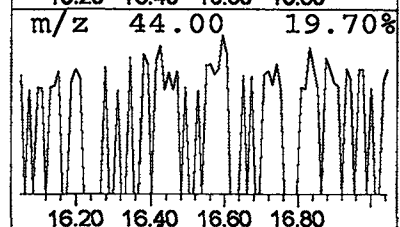
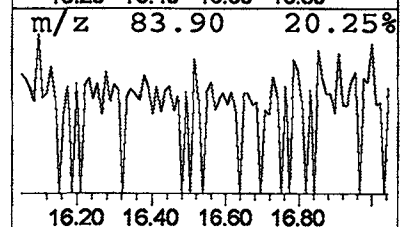
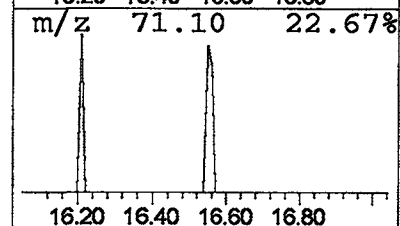
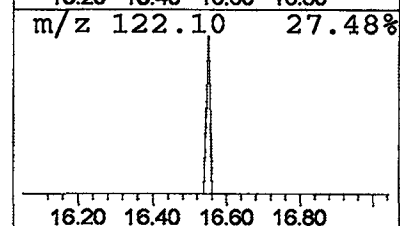
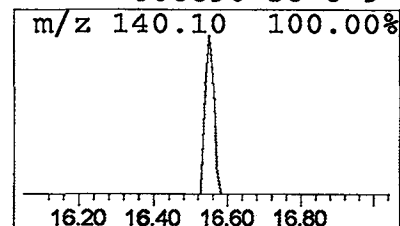
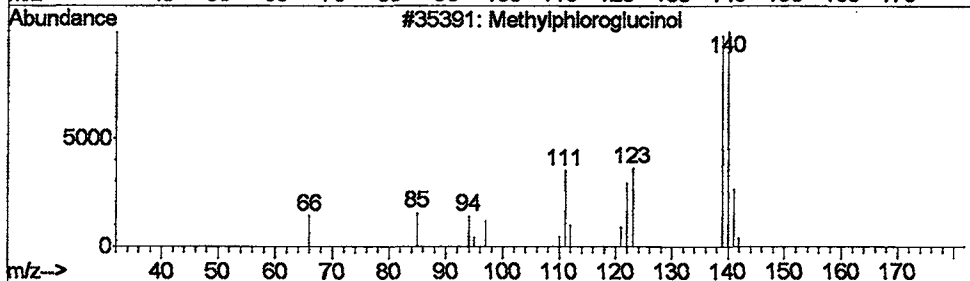
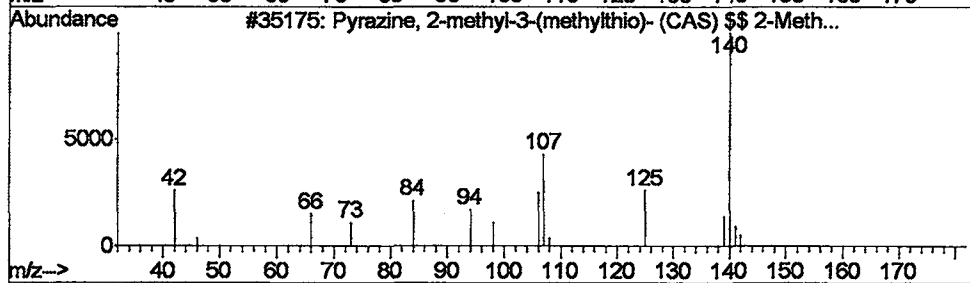
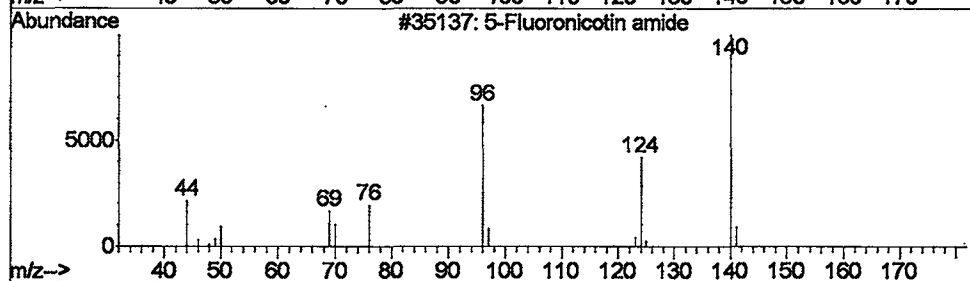
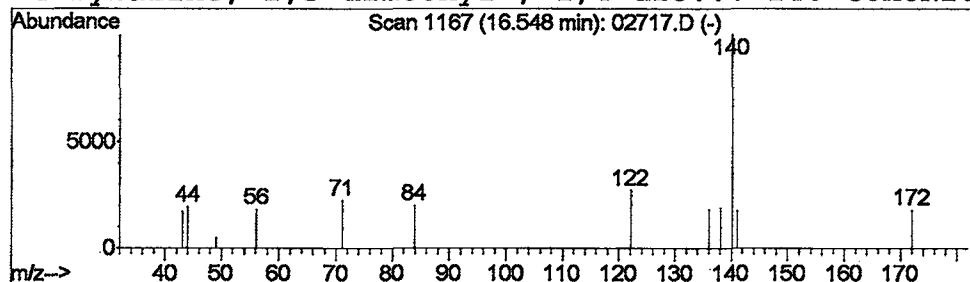
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 Inst : Instrumen
 Multiplr: 1.00

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

 Peak Number 10 5-Fluoronicotin amide Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.55	0.03 ug/L	18866	Acenaphthene-d10	18.33

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	5-Fluoronicotin amide	140	C6H5FN2O	000000-00-0	25
2	Pyrazine, 2-methyl-3-(methylthio...	140	C6H8N2S	002882-20-4	9
3	Methylphloroglucinol	140	C7H8O3	000000-00-0	9
4	Pyrazine, 2,5-dimethyl-, 1,4-dio...	140	C6H8N2O2	006890-38-6	9



Library Search Compound Report

Data File : C:\MSDCHEM\1\5973N\02FEB28\02717.D
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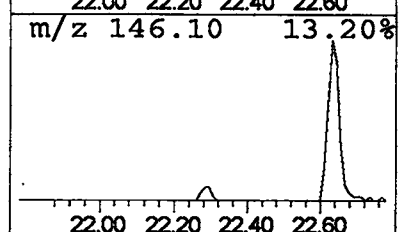
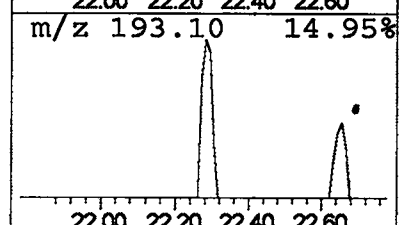
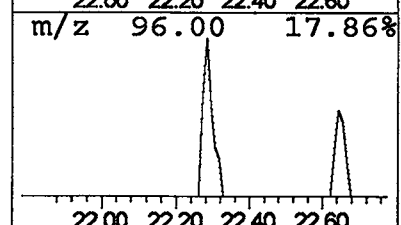
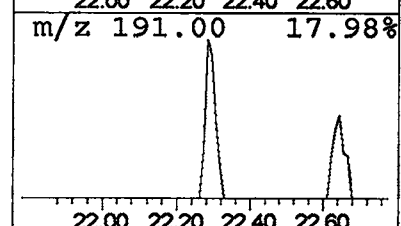
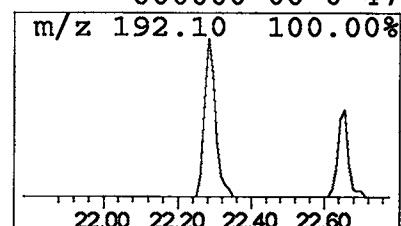
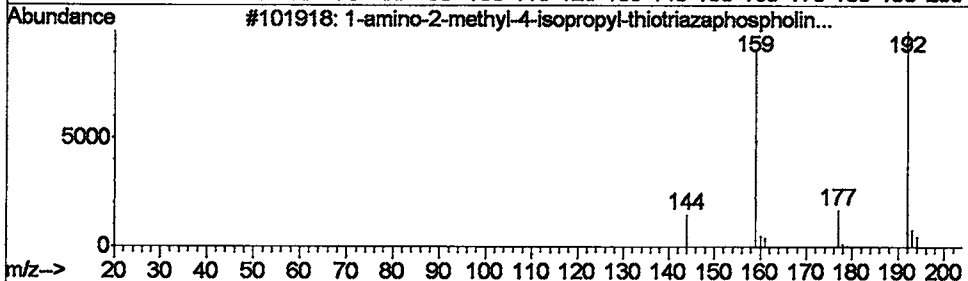
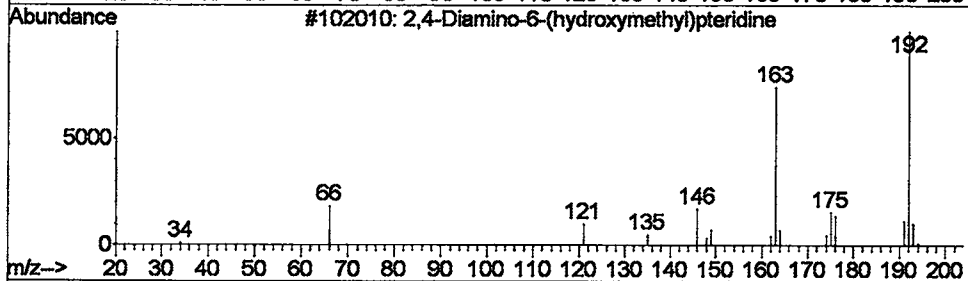
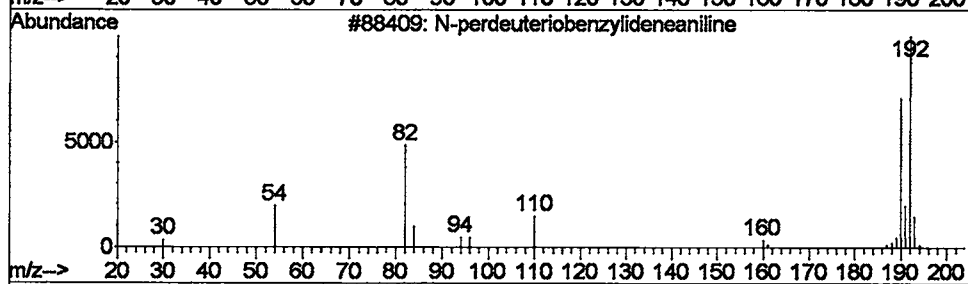
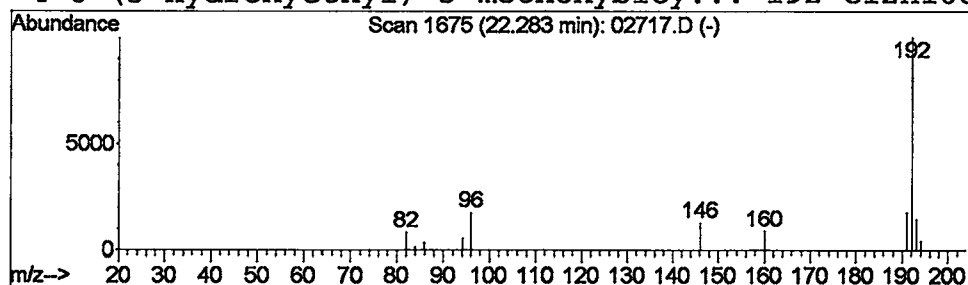
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 Multiplr: 1.00

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

 Peak Number 11 N-perdeuteriobenzylideneani... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.28	0.10 ug/L	62034	Phenanthrene-d10	22.63

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		N-perdeuteriobenzylideneaniline	192	C13D11N	000538-51-2	64
2		2,4-Diamino-6-(hydroxymethyl)pte...	192	C7H8N6O	000945-24-4	50
3		1-amino-2-methyl-4-isopropyl-thi...	192	C5H13N4PS	107264-54-0	50
4		8-(3-Hydroxyethyl)-3-methoxybicy...	192	C12H16O2	000000-00-0	47



Library Search Compound Report

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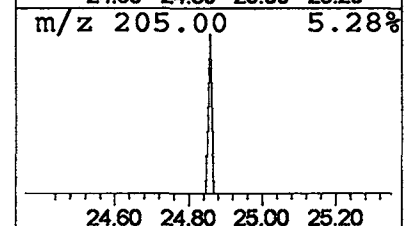
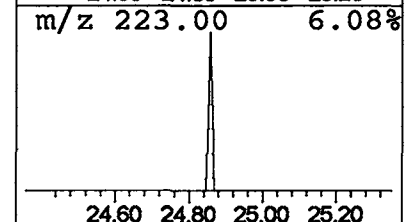
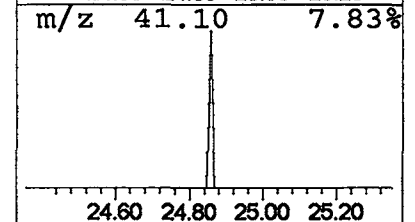
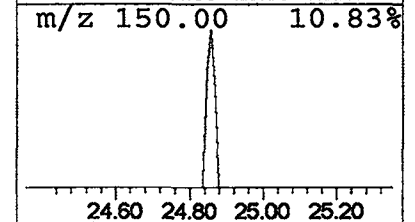
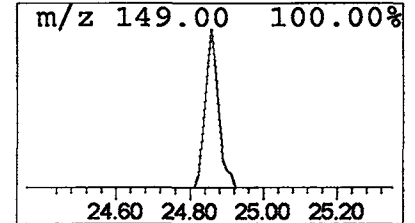
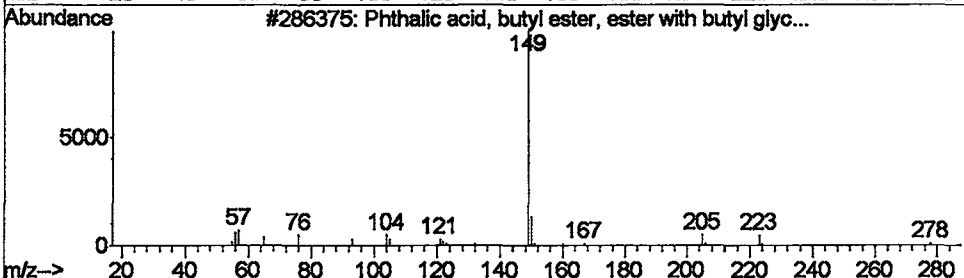
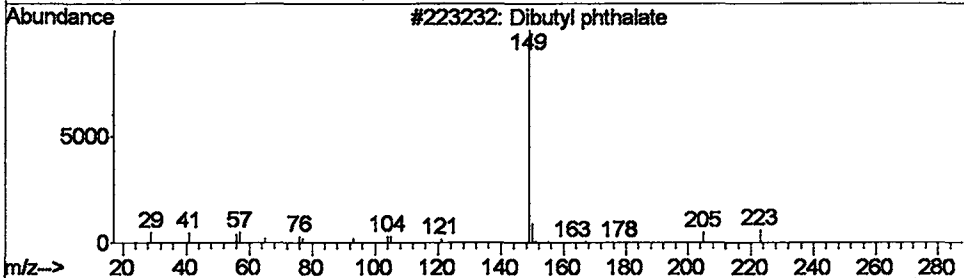
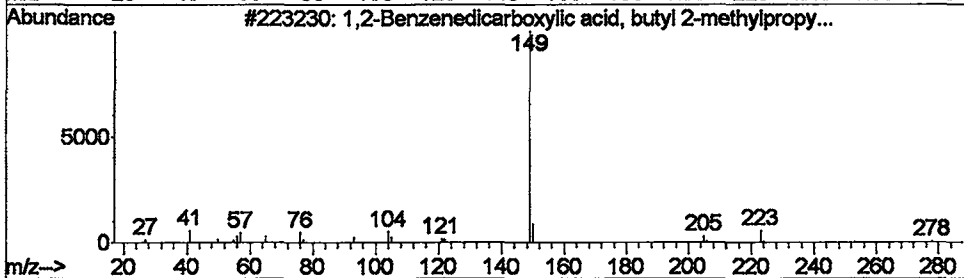
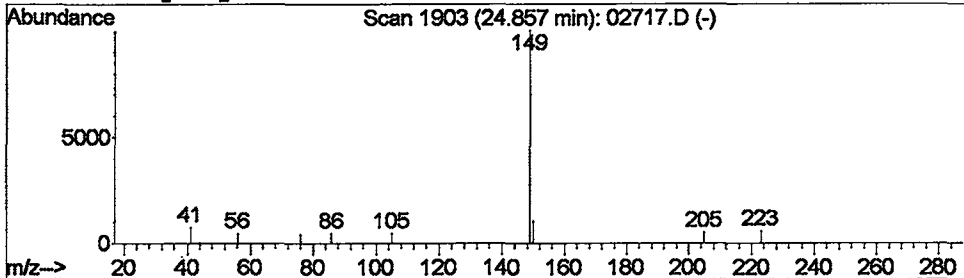
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Quant Method : C:\MSDCHEM\1...\HP022102.M (RTE Integrator)
 Title : 508 scan method
 Library : C:\DATABASE\WILEY7N.L

 Peak Number 12 1,2-Benzenedicarboxylic aci... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.86	0.06 ug/L	38545	Phenanthrene-d10	22.63

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		1,2-Benzenedicarboxylic acid, bu...	278	C16H22O4	017851-53-5	83
2		Dibutyl phthalate	278	C16H22O4	000084-74-2	83
3		Phthalic acid, butyl ester, este...	336	C18H24O6	000085-70-1	83
4		Dibutyl phthalate	278	C16H22O4	000084-74-2	74



Library Search Compound Report

Data File : C:\MSDCHEM\1\5973N\02FEB28\02717.D
 Acq On : 28 Feb 2002 3:30 pm
 Sample : 508 scan
 Misc : February 28, 2002
 MS Integration Params: LSCINT.P

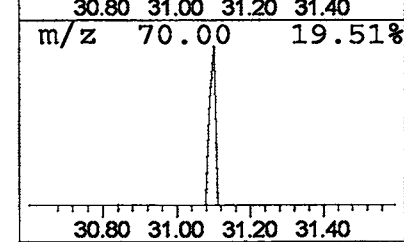
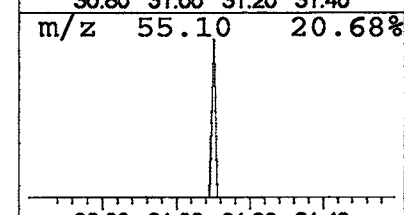
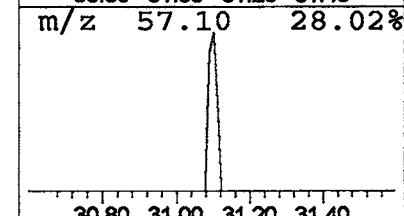
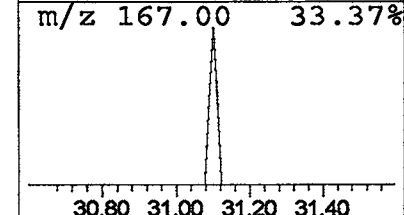
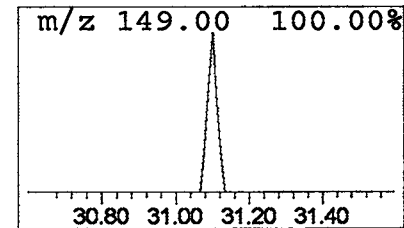
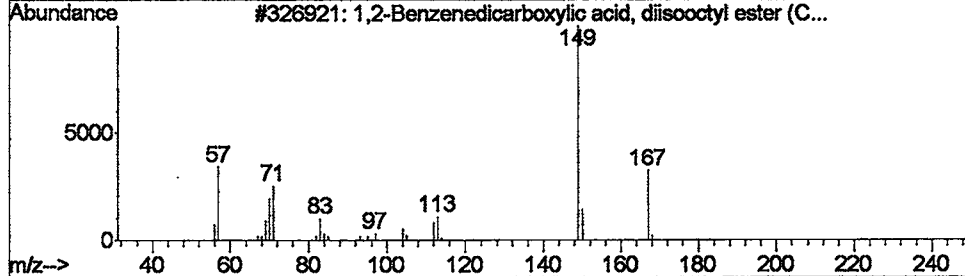
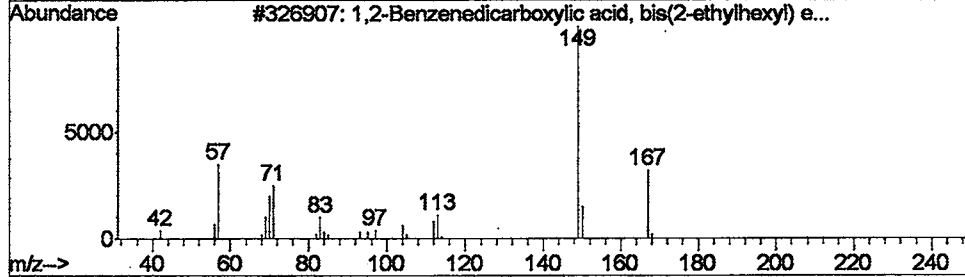
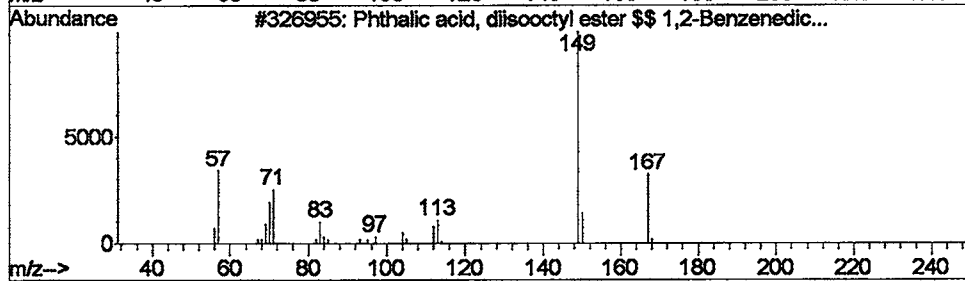
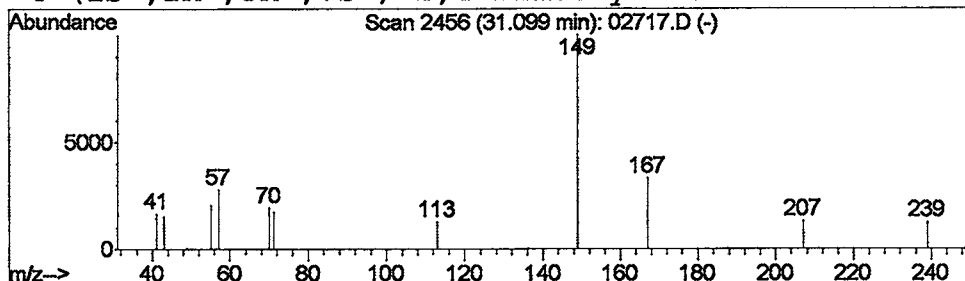
Vial: 2
 Operator:
 Inst : Instrumen
 Multiplr: 1.00

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

 Peak Number 13 Phthalic acid, diisooctyl e... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
31.10	0.04 ug/L	23756	Chrysene-d12	30.49

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Phthalic acid, diisooctyl ester ...	390	C24H38O4	001330-91-2	72
2		1,2-Benzenedicarboxylic acid, bi...	390	C24H38O4	000117-81-7	72
3		1,2-Benzenedicarboxylic acid, di...	390	C24H38O4	027554-26-3	72
4		(1S*,2R*,5R*,7S*)-2,4-Dimethyl-7...	168	C10H16O2	091926-28-2	64



Tentatively Identified Compound (LSC) summary

Operator ID: Date Acquired: 28 Feb 2002 3:30 pm
 Data File: C:\MSDCHEM\1\5973N\02FEB28\02717.D
 Name: 508 scan
 Misc: February 28, 2002
 Method: C:\MSDCHEM\1\METHODS\METHODS\HP022102.M (RTE Integrator)
 Title: 508 scan method
 Library Searched: C:\DATABASE\WILEY7N.L

TIC Top Hit name	RT	EstConc	Units	Response	#	--Internal Standard---		
						RT	Resp	Conc
Butanoic acid, me...	3.61	0.2	ug/L	56700	1	9.72	7318690	20.0
ISOBUTYRALDEHYDE,...	3.92	0.1	ug/L	27400	1	9.72	7318690	20.0
Cyclopropanecarbo...	3.99	0.0	ug/L	17400	1	9.72	7318690	20.0
2-Butenoic acid, ...	4.24	0.0	ug/L	17474	1	9.72	7318690	20.0
Acetic acid, 3-[1...	5.94	0.3	ug/L	126559	1	9.72	7318690	20.0
CH3C(O)NHCO(O)CH3 ...	6.12	0.1	ug/L	29490	1	9.72	7318690	20.0
1-DEUTEROBUTANE	9.89	0.1	ug/L	30498	1	9.72	7318690	20.0
Cyclopentasiloxan...	12.54	0.0	ug/L	17488	2	13.20	10476700	20.0
Pyrazine, 2,5-dim...	13.38	0.0	ug/L	20438	2	13.20	10476700	20.0
5-Fluoronicotin a...	16.55	0.0	ug/L	18866	3	18.33	11175700	20.0
N-perdeuteriobenz...	22.28	0.1	ug/L	62034	4	22.63	11949600	20.0
1,2-Benzenedicarb...	24.86	0.1	ug/L	38545	4	22.63	11949600	20.0
Phthalic acid, di...	31.10	0.0	ug/L	23756	5	30.49	11147200	20.0