

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

Sample: AE02716
 Analysis: \$GCMS GCMS Scan For Unknowns
 Units: ug
 Started: 3/5/02 12:02 Ended: 3/5/02 12:02 Analyst: DLV
 Page: 1

	Component Name	Viol	Result	Qualifier	MDL
1	Other unknown compounds		.		
2	Surr_2,4-DDD		102		5.0

Comments for Sample AE02716 - Analysis \$GCMS

Westchester Dept. of Labs & Research

User: Vinci, David L

Date: 03-05-2002 Time: 12:04:35

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\02716.D Vial: 1
 Acq Cn : 28 Feb 2002 2:40 pm Operator:
 Sample : 508 scan Inst : Instrumen
 Misc : February 28, 2002 Multiplr: 1.00
 MS Integration Params: SPLIT.P
 Quant Time: Mar 01 09:02:54 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	1312634	20.00	ug/L	0.00
10) Naphthalene-d8	13.20	136	5739558	20.00	ug/L	0.00
17) Acenaphthene-d10	18.33	164	2902133	20.00	ug/L	0.00
29) Phenanthrene-d10	22.63	188	4883853	20.00	ug/L	-0.01
38) Chrysene-d12	30.49	240	3904088	20.00	ug/L	-0.03
44) Perylene-d12	34.41	264	2486483	20.00	ug/L	-0.04

System Monitoring Compounds
 37) 2,4-DDD 27.72 235 333193 5.12 ug/L 0.00
 Spiked Amount 5.000 Recovery = 102.40%

Target Compounds
 20) Dimethylphthalate 18.34 163 468304 2.66 ug/L # 57
 21) 2,6-Dinitrotoluene 18.33 165 362843 8.92 ug/L # 27

Quantitation Report

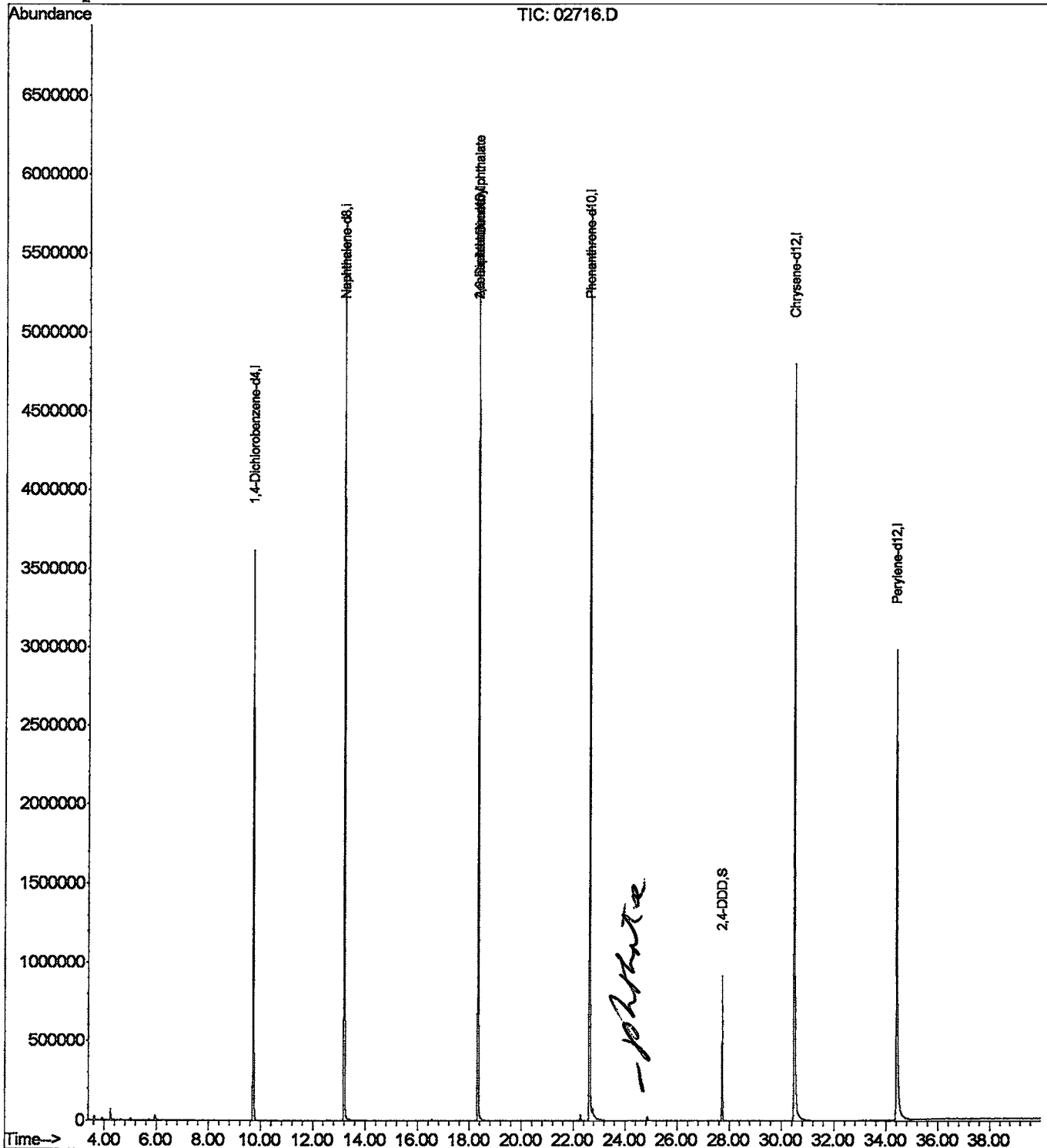
(Not Reviewed)

Data File : C:\MSDCHEM\1\5973N\02FEB28\02716.D
 Acq On : 28 Feb 2002 2:40 pm
 Sample : 508 scan
 Misc : February 28, 2002
 MS Integration Params: SPLIT.P
 Quant Time: Mar 1 9:02 2002

Vial: 1
 Operator:
 Inst : Instrumen
 Multiplr: 1.00

Quant Results File: HP022102.R

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



LSC Area Percent Report

Data File : C:\MSDCHEM\1\5973N\02FEB28\02716.D
 Acq On : 28 Feb 2002 2:40 pm
 Sample : 508 scan
 Misc : February 28, 2002
 MS Integration Params: LSCINT.P

Vial: 1
 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

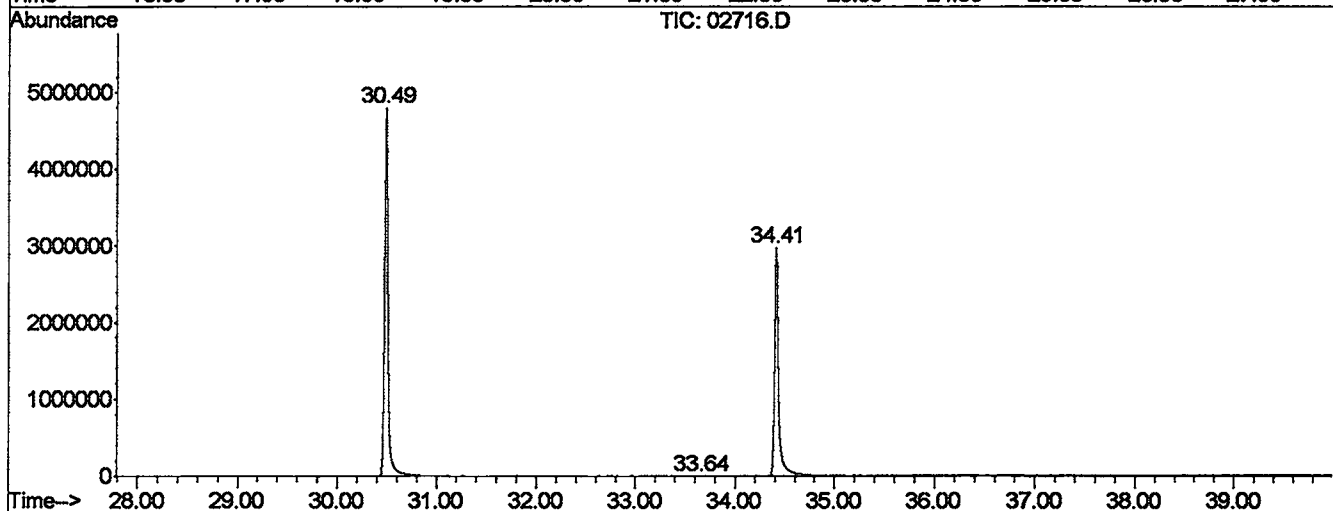
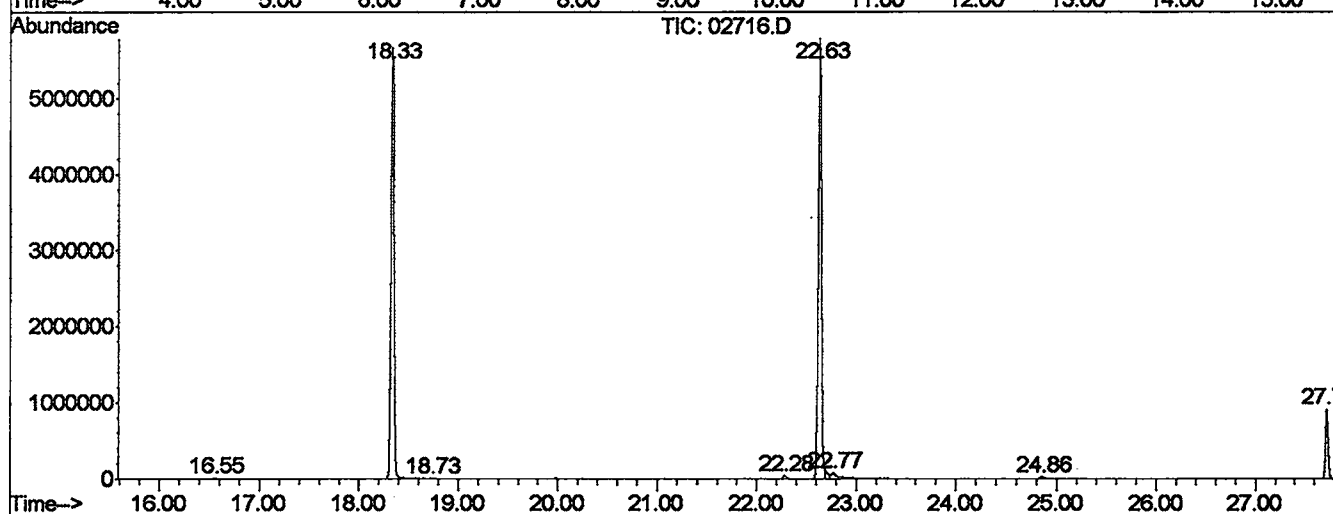
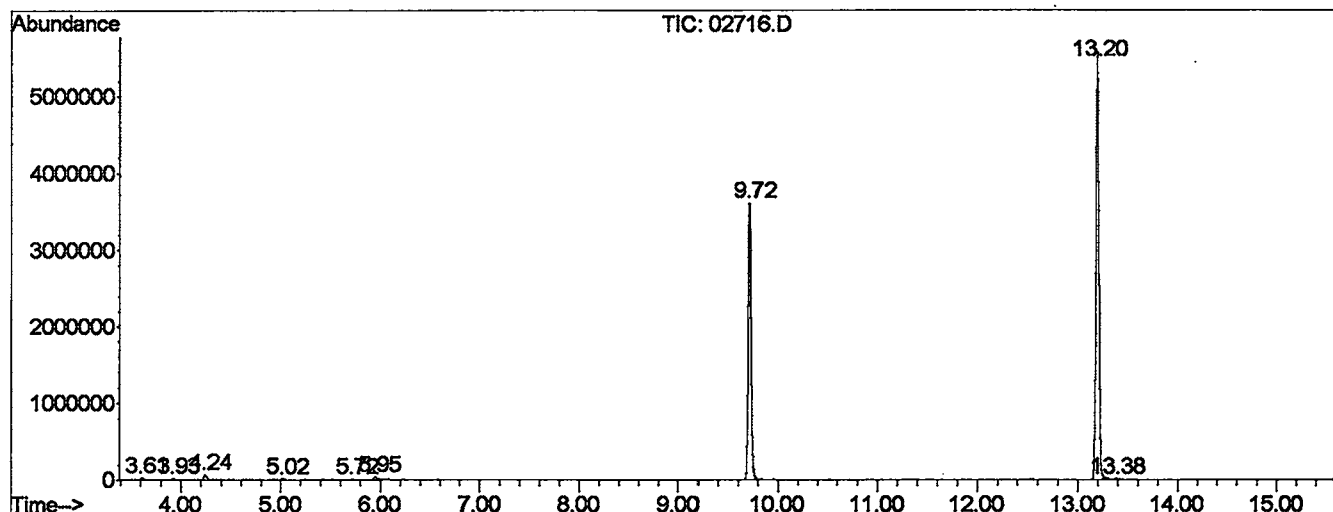
Signal : TIC

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	31	rVB	26626	50459	0.40%	0.077%
2	3.927	45	49	53	rBV	17004	28562	0.23%	0.044%
3	4.243	73	77	85	rVV3	69596	128314	1.01%	0.197%
4	5.022	143	146	151	rBV3	13502	29102	0.23%	0.045%
5	5.722	205	208	213	rVB4	7244	13724	0.11%	0.021%
6	5.948	223	228	249	rVB	32715	107145	0.84%	0.164%
7	9.718	557	562	567	rBV	3612595	7355022	57.94%	11.279%
8	13.195	865	870	875	rBV	5577545	10780971	84.93%	16.533%
9	13.376	883	886	891	rVB2	10002	25450	0.20%	0.039%
10	16.548	1163	1167	1173	rBV3	9604	21076	0.17%	0.032%
11	18.332	1319	1325	1331	rBV	5672087	11831537	93.20%	18.144%
12	18.727	1357	1360	1367	rVB4	4414	15611	0.12%	0.024%
13	22.283	1669	1675	1679	rBV	37706	75603	0.60%	0.116%
14	22.633	1701	1706	1711	rBV2	5784305	12694180	100.00%	19.467%
15	22.768	1715	1718	1751	rVB	74486	396144	3.12%	0.607%
16	24.857	1899	1903	1907	rBV	22991	47657	0.38%	0.073%
17	27.724	2153	2157	2163	rBV	911885	1601364	12.61%	2.456%
18	30.490	2397	2402	2441	rBV	4794175	11517042	90.73%	17.661%
19	33.639	2671	2681	2683	rBV4	3155	17248	0.14%	0.026%
20	34.407	2743	2749	2783	rBV	2975757	8473812	66.75%	12.995%

Sum of corrected areas: 65210023

LSC Report - Integrated Chromatogram

File : C:\MSDCHEM\1\5973N\02FEB28\02716.D
 Operator :
 Acquired : 28 Feb 2002 2:40 pm using AcqMethod HP020702
 Instrument : Instrumen
 Sample Name: 508 scan
 Misc Info : February 28, 2002
 Vial Number: 1
 Quant File :HP022102.RES (RTE Integrator)



Library Search Compound Report

Data File : C:\MSDCHEM\1\5973N\02FEB28\02716.D
 Acq On : 28 Feb 2002 2:40 pm
 Sample : 508 scan
 Misc : February 28, 2002
 MS Integration Params: LSCINT.P

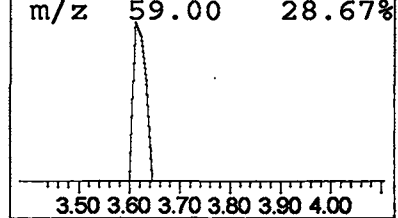
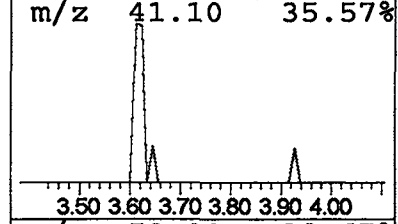
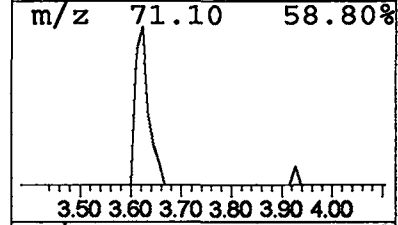
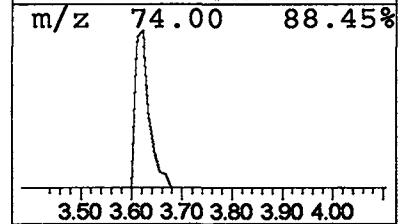
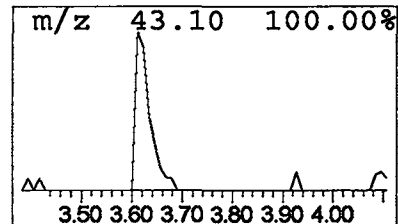
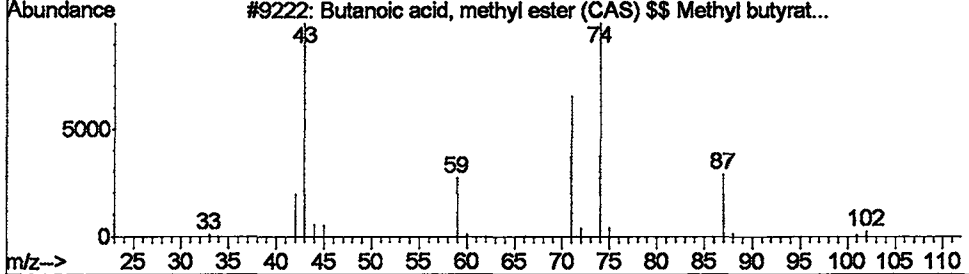
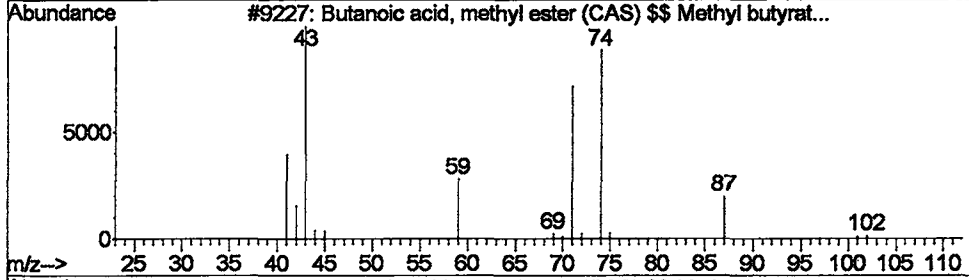
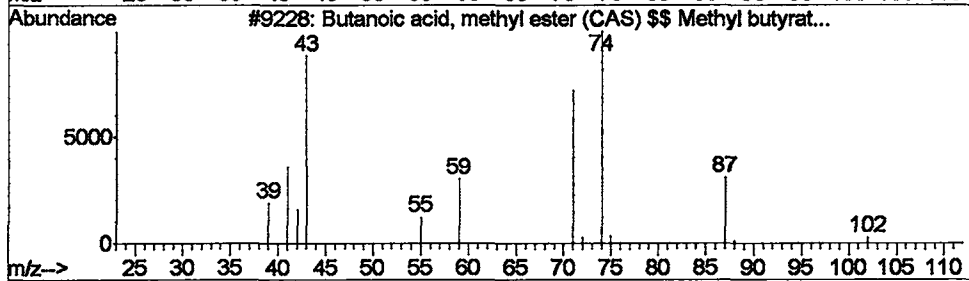
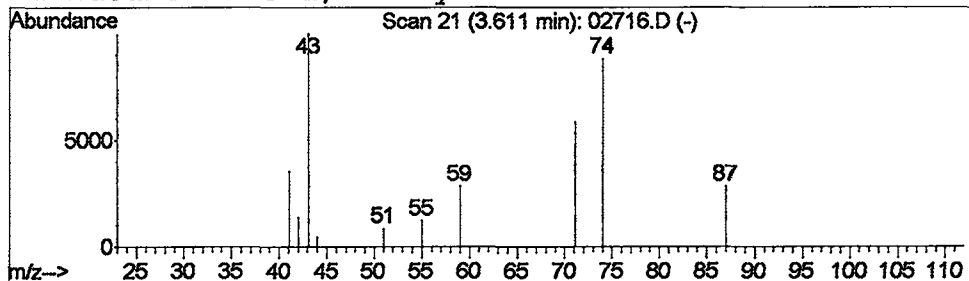
Vial: 1
 Operator:
 Inst : Instrumen
 Multiplr: 1.00

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 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.61	0.14 ug/L	50459	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 (CAS...	102	C5H10O2	000623-42-7	83
4			Butanoic acid, methyl ester	102	C5H10O2	000623-42-7	78



Library Search Compound Report

Data File : C:\MSDCHEM\1\5973N\02FEB28\02716.D
 Acq On : 28 Feb 2002 2:40 pm
 Sample : 508 scan
 Misc : February 28, 2002
 MS Integration Params: LSCINT.P

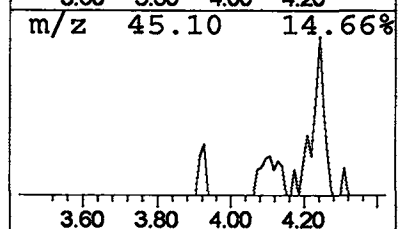
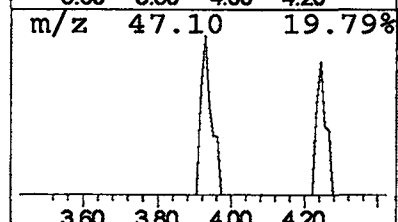
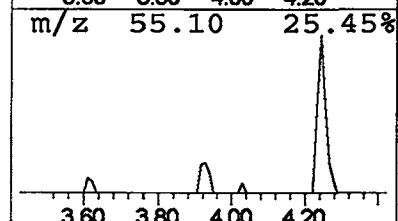
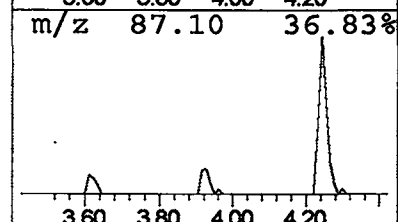
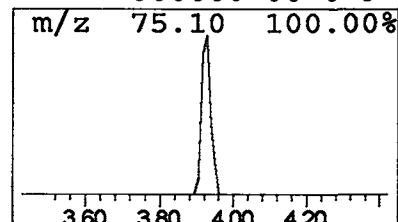
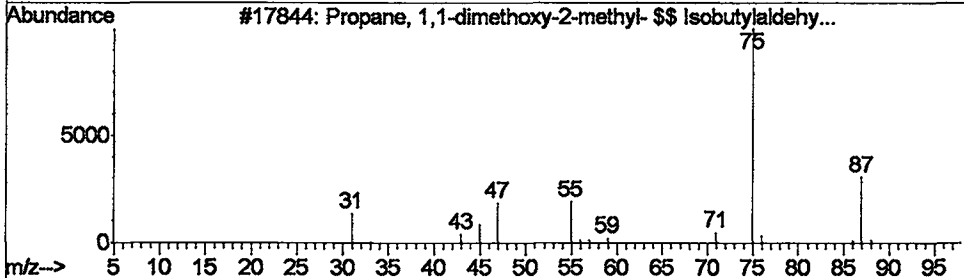
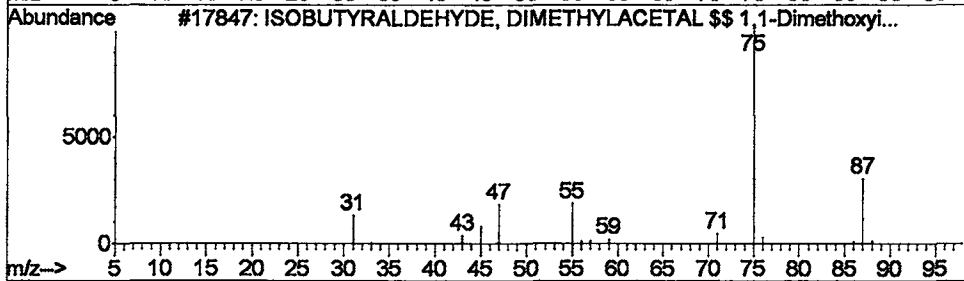
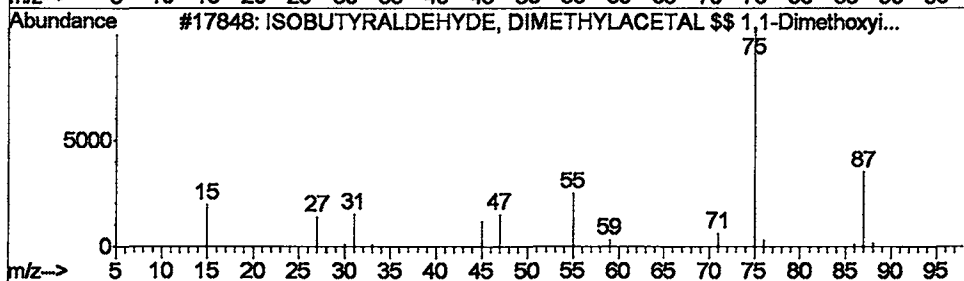
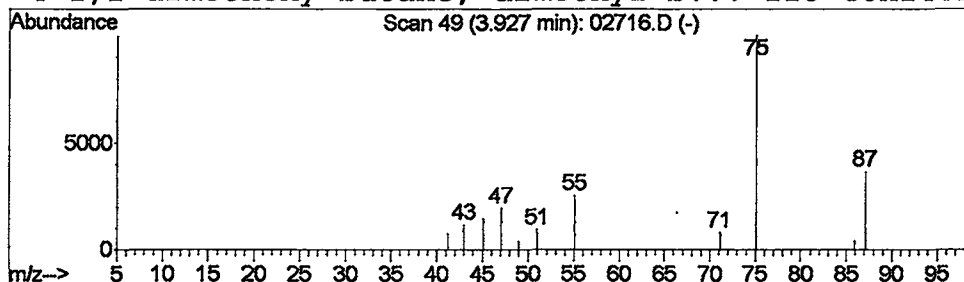
Vial: 1
 Operator:
 Inst : Instrumen
 Multiplr: 1.00

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.93	0.08 ug/L	28562	1,4-Dichlorobenzene-d4	9.72

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		ISOBUTYRALDEHYDE, DIMETHYLACETAL...	118	C6H14O2	000000-00-0	72
2		ISOBUTYRALDEHYDE, DIMETHYLACETAL...	118	C6H14O2	000000-00-0	56
3		Propane, 1,1-dimethoxy-2-methyl...	118	C6H14O2	041632-89-7	56
4		1,1-dimethoxy butane, dimethyl b...	118	C6H14O2	000000-00-0	9



Library Search Compound Report

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 Acq On : 28 Feb 2002 2:40 pm
 Sample : 508 scan
 Misc : February 28, 2002
 MS Integration Params: LSCINT.P

Vial: 1
 Operator:
 Inst : Instrumen
 Multiplr: 1.00

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

Title : 508 scan method

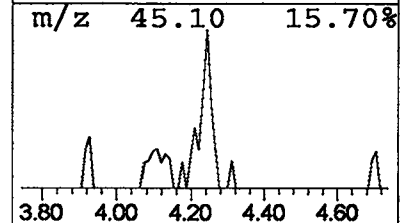
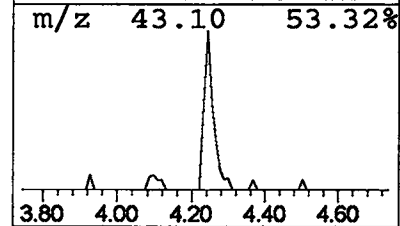
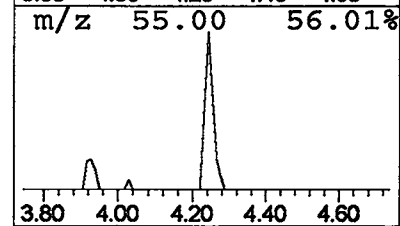
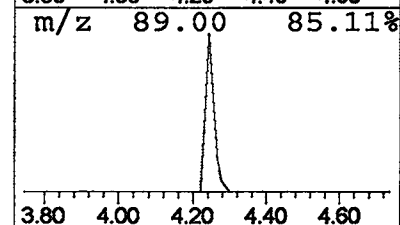
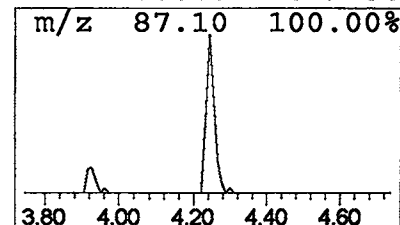
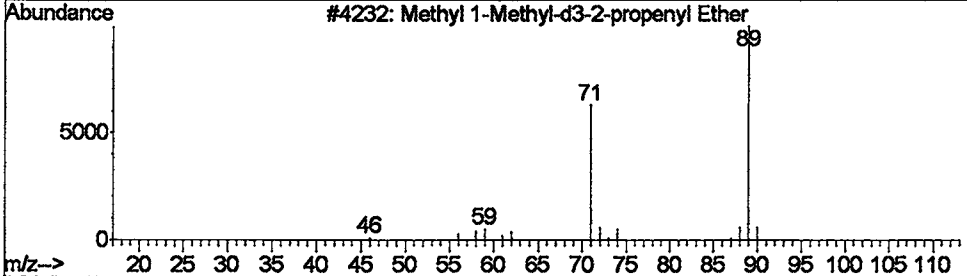
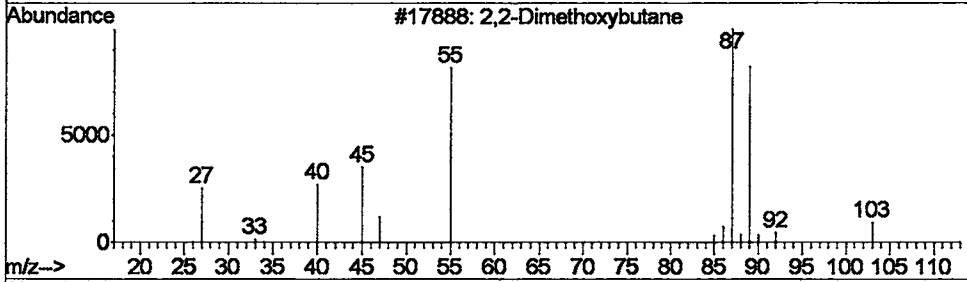
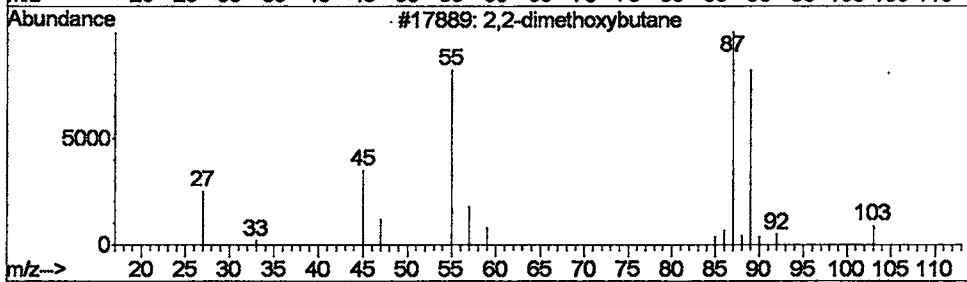
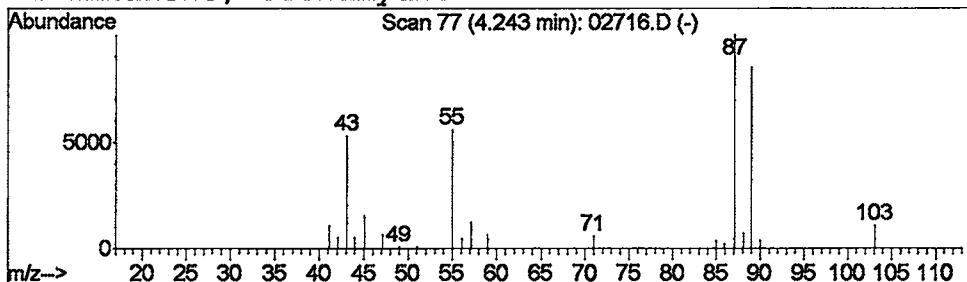
Library : C:\DATABASE\WILEY7N.L

 Peak Number 3 2,2-dimethoxybutane

Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.24	0.35 ug/L	128314	1,4-Dichlorobenzene-d4	9.72

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2,2-dimethoxybutane	118	C6H14O2	003453-99-4	83
2			2,2-Dimethoxybutane	118	C6H14O2	003453-99-4	74
3			Methyl 1-Methyl-d3-2-propenyl Ether	89	C5H7D3O	000000-00-0	42
4			Thiazole, tetrahydro-	89	C3H7NS	000504-78-9	36



Library Search Compound Report

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

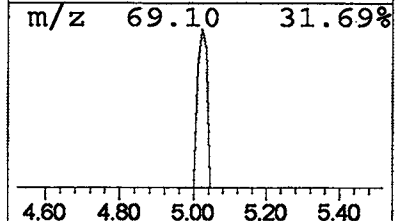
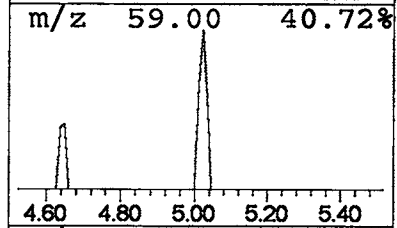
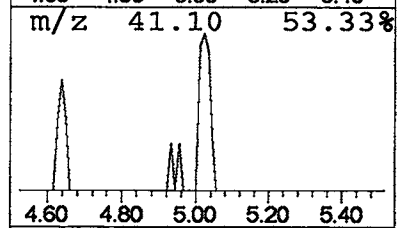
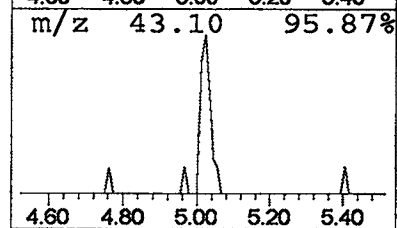
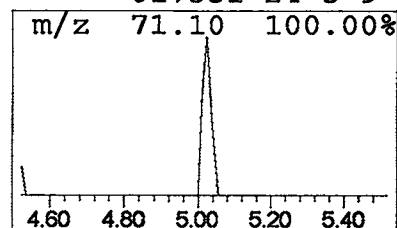
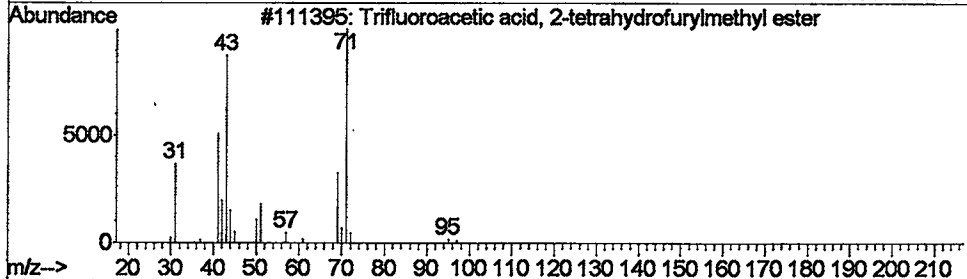
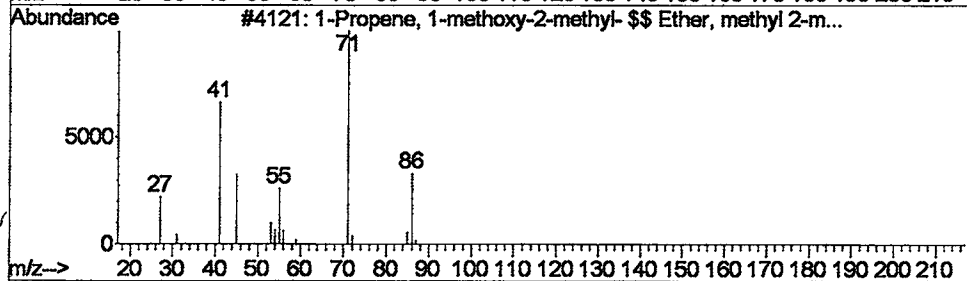
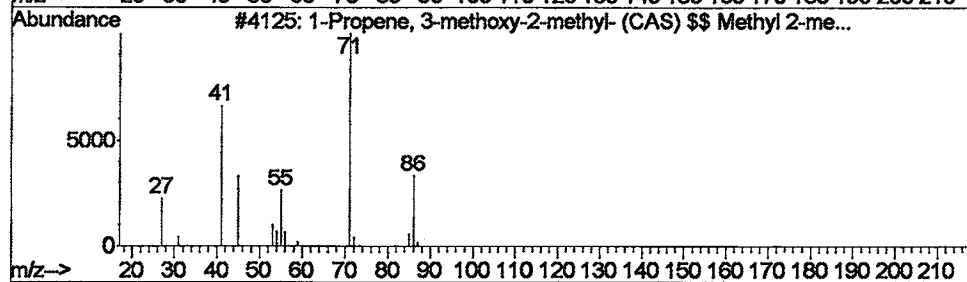
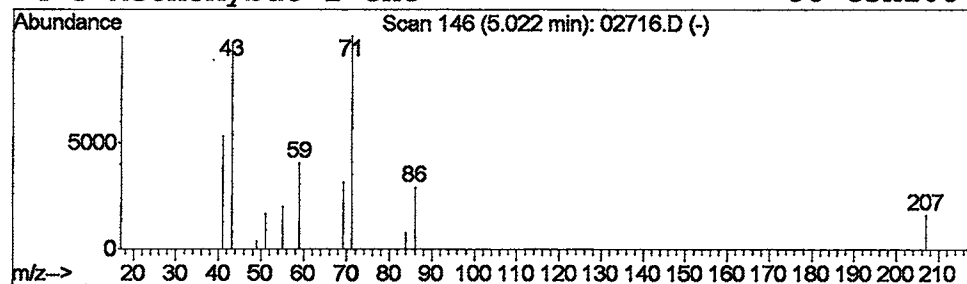
Vial: 1
 Operator:
 Inst : Instrumen
 Multiplr: 1.00

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

 Peak Number 4 1-Propene, 3-methoxy-2-meth... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.02	0.08 ug/L	29102	1,4-Dichlorobenzene-d4	9.72

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1-Propene, 3-methoxy-2-methyl- (...	86	C5H10O	022418-49-1	38
2			1-Propene, 1-methoxy-2-methyl- \$...	86	C5H10O	017574-84-4	38
3			Trifluoroacetic acid, 2-tetrahyd...	198	C7H9F3O3	000000-00-0	23
4			3-Methoxybut-1-ene	86	C5H10O	017351-24-5	9



Library Search Compound Report

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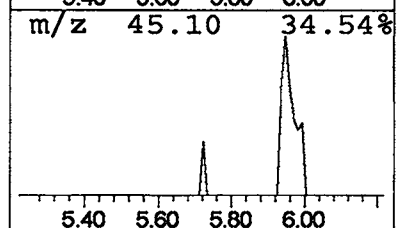
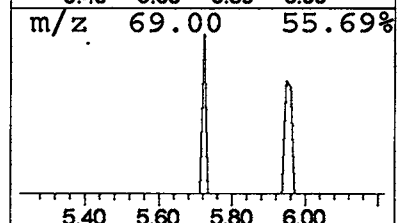
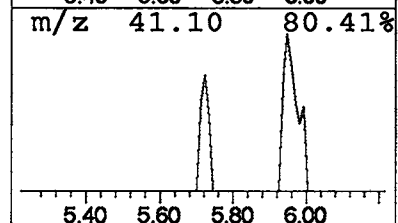
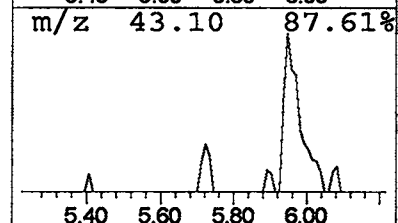
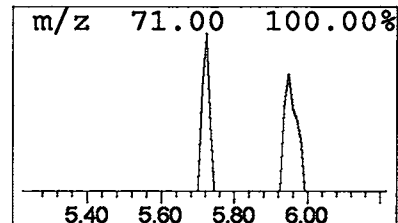
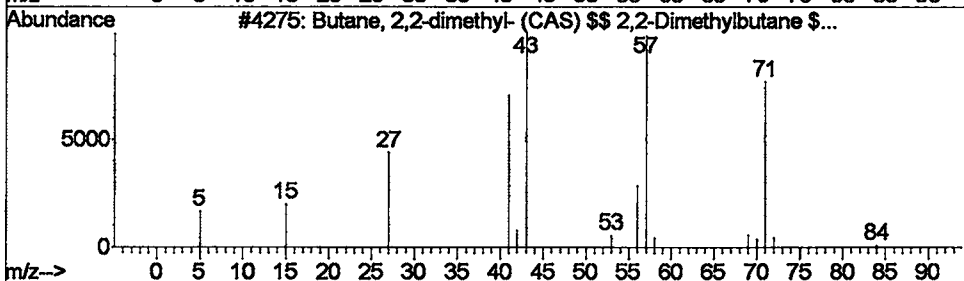
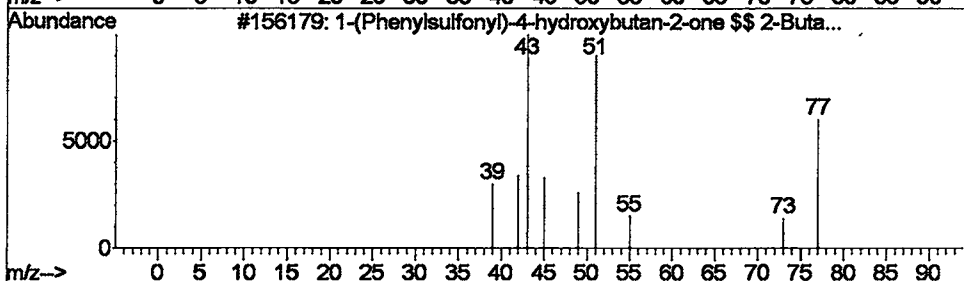
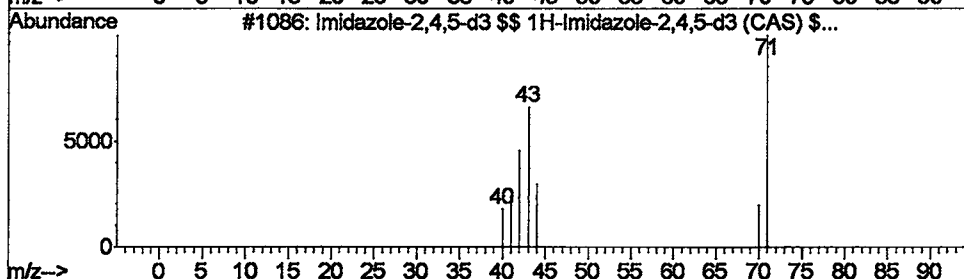
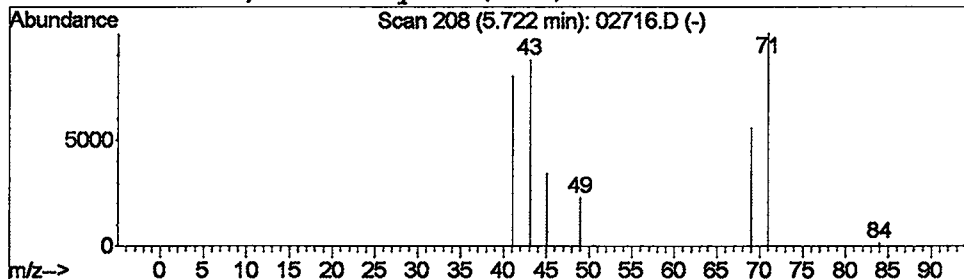
Vial: 1
Operator:
Inst : Instrumen
Multiplr: 1.00

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

Peak Number 5 Imidazole-2,4,5-d3 \$\$ 1H-Im... Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.72	0.04 ug/L	13724	1,4-Dichlorobenzene-d4	9.72

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Imidazole-2,4,5-d3 \$\$ 1H-Imidazo...	71	C3HD3N2	006745-43-3	4
2			1-(Phenylsulfonyl)-4-hydroxybuta...	228	C10H12O4S	088703-48-4	4
3			Butane, 2,2-dimethyl- (CAS) \$\$ 2...	86	C6H14	000075-83-2	4
4			Azetidine, 2-methyl- (CAS)	71	C4H9N	019812-49-8	3



Library Search Compound Report

Data File : C:\MSDCHEM\1\5973N\02FEB28\02716.D
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 Sample : 508 scan
 Misc : February 28, 2002
 MS Integration Params: LSCINT.P

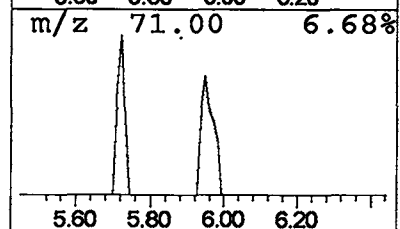
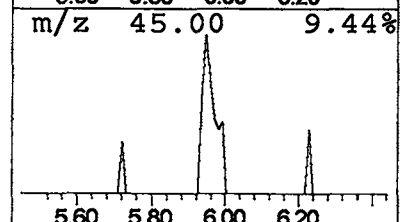
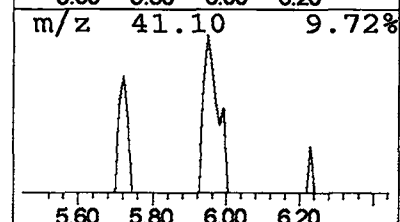
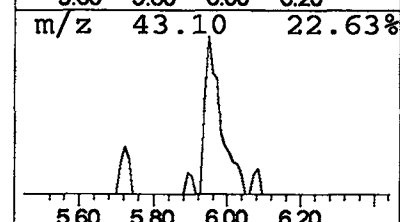
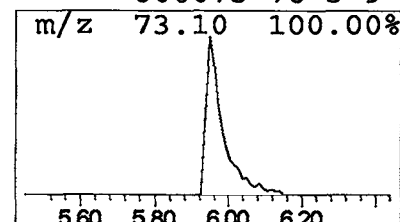
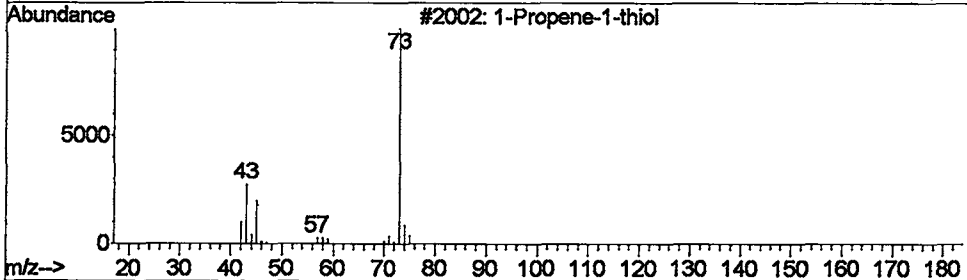
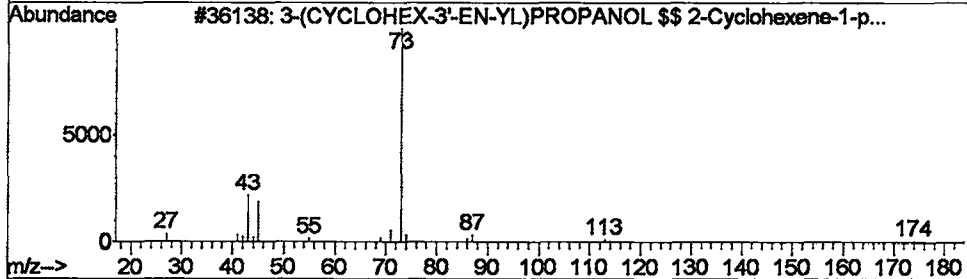
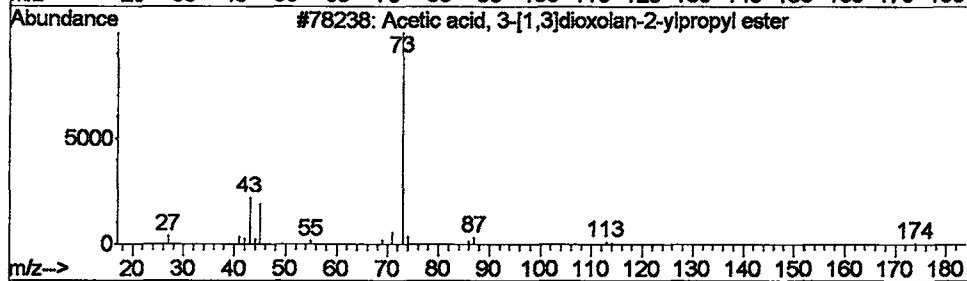
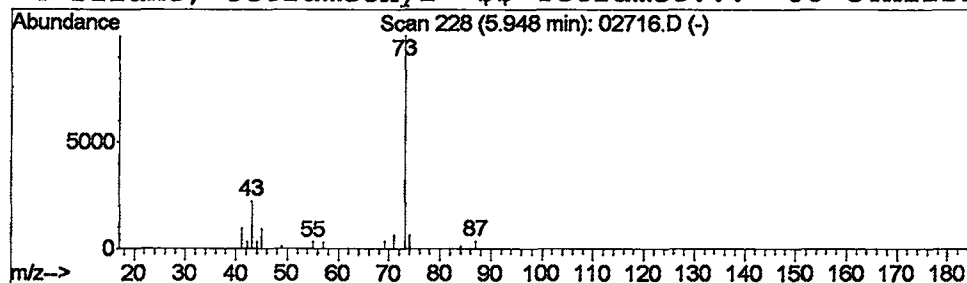
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Quant Method : C:\MSDCHEM\1...\HP022102.M (RTE Integrator)
 Title : 508 scan method
 Library : C:\DATABASE\WILEY7N.L

 Peak Number 6 Acetic acid, 3-[1,3]dioxola... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.95	0.29 ug/L	107145	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	38
2	3-(CYCLOHEX-3'-EN-YL)PROPANOL \$\$...	174	C8H14O4	015745-87-6	38
3	1-Propene-1-thiol	74	C3H6S	000925-89-3	9
4	Silane, tetramethyl- \$\$ Tetramet...	88	C4H12Si	000075-76-3	9



Library Search Compound Report

Data File : C:\MSDCHEM\1\5973N\02FEB28\02716.D
 Acq On : 28 Feb 2002 2:40 pm
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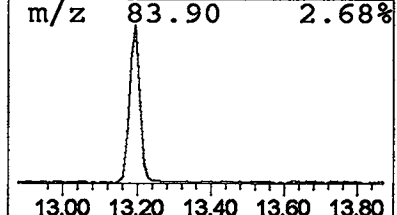
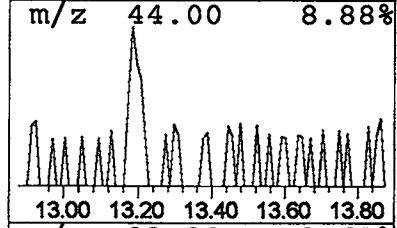
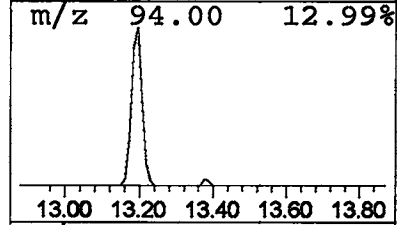
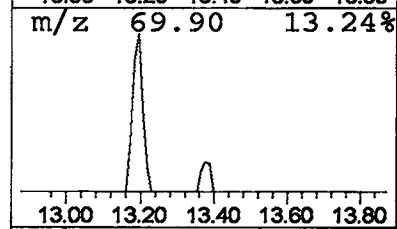
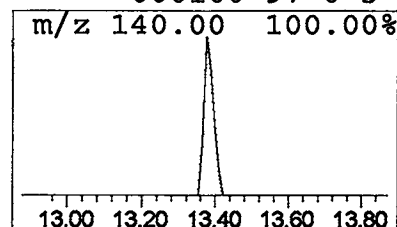
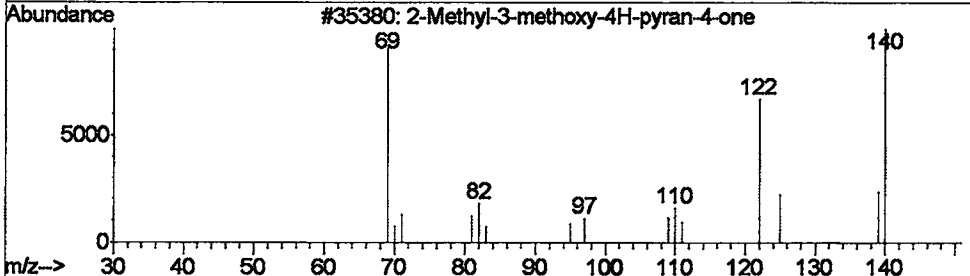
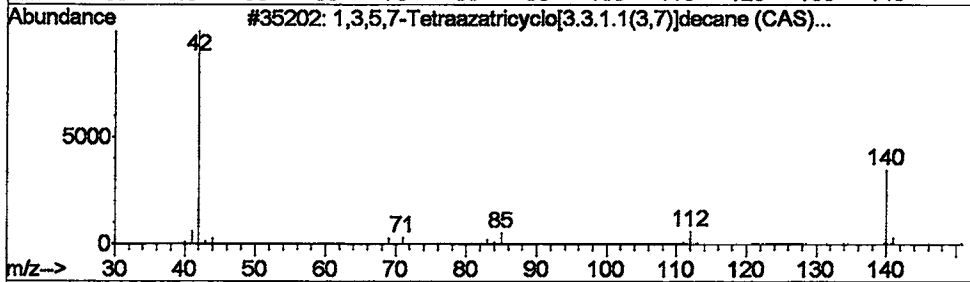
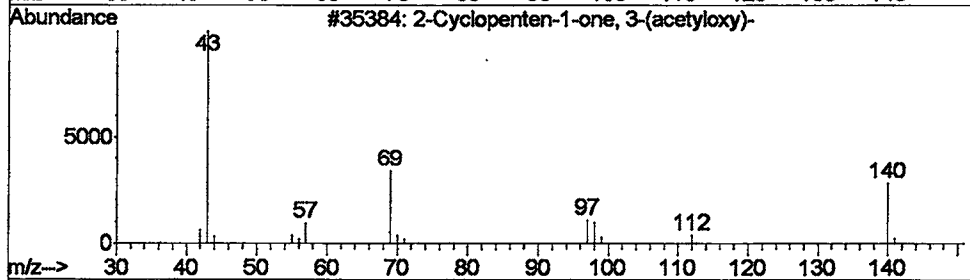
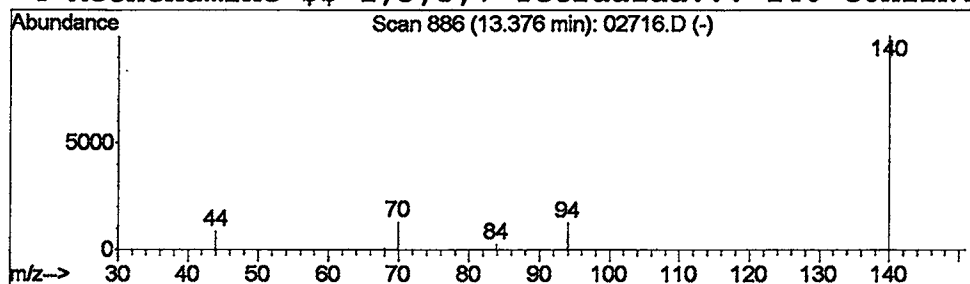
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Quant Method : C:\MSDCHEM\1...\HP022102.M (RTE Integrator)
 Title : 508 scan method
 Library : C:\DATABASE\WILEY7N.L

 Peak Number 7 2-Cyclopenten-1-one, 3-(ace... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.38	0.05 ug/L	25450	Naphthalene-d8	13.20

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Cyclopenten-1-one, 3-(acetyloxy)-	140	C7H8O3	018766-96-6	5
2		1,3,5,7-Tetraazatricyclo[3.3.1.1...]	140	C6H12N4	000100-97-0	4
3		2-Methyl-3-methoxy-4H-pyran-4-one	140	C7H8O3	000000-00-0	4
4		Methenamine \$\$ 1,3,5,7-Tetraazaa...	140	C6H12N4	000100-97-0	3



Library Search Compound Report

Data File : C:\MSDCHEM\1\5973N\02FEB28\02716.D
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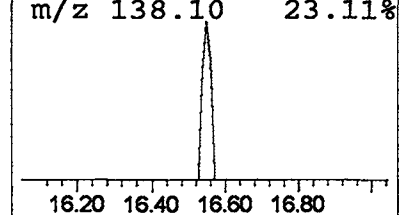
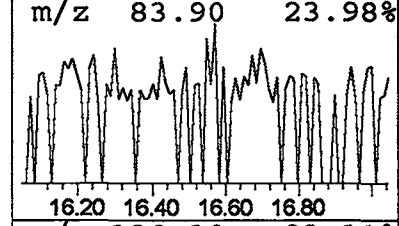
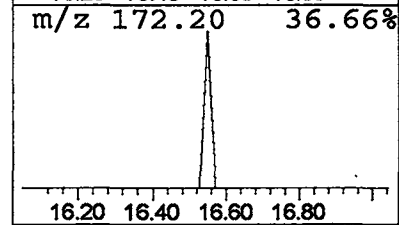
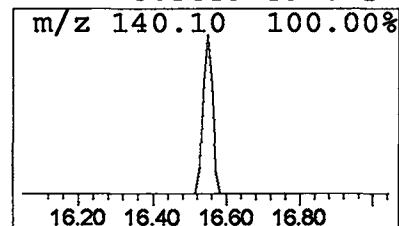
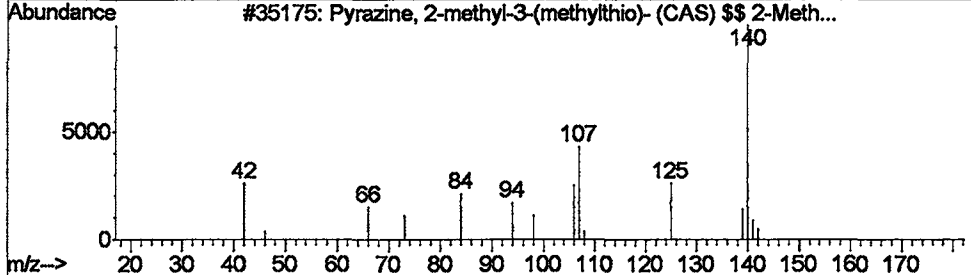
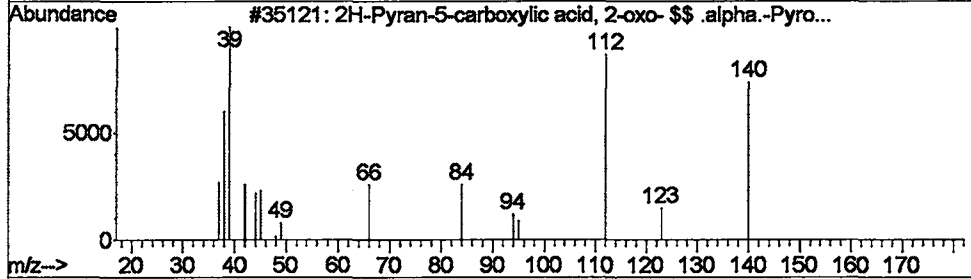
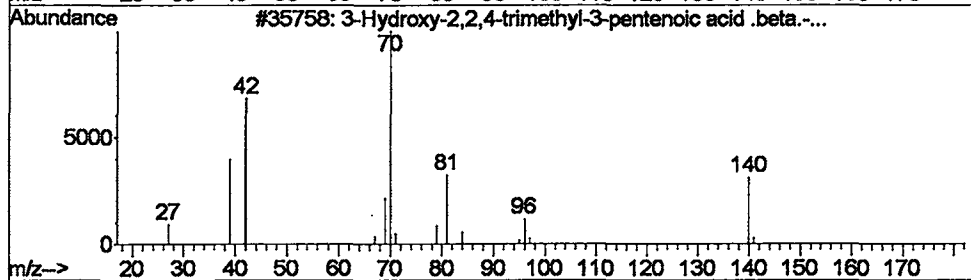
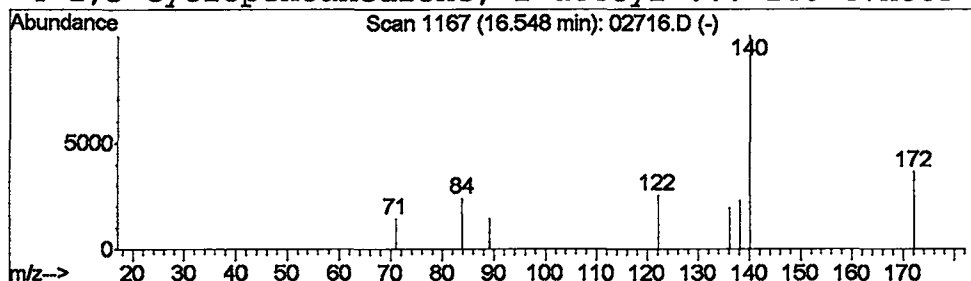
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Quant Method : C:\MSDCHEM\1...\HP022102.M (RTE Integrator)
 Title : 508 scan method
 Library : C:\DATABASE\WILEY7N.L

 Peak Number 8 3-Hydroxy-2,2,4-trimethyl-3... Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.55	0.04 ug/L	21076	Acenaphthene-d10	18.33

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Hydroxy-2,2,4-trimethyl-3-pent...	140	C8H12O2	003173-79-3	5
2		2H-Pyran-5-carboxylic acid, 2-ox...	140	C6H4O4	000500-05-0	5
3		Pyrazine, 2-methyl-3-(methylthio...	140	C6H8N2S	002882-20-4	5
4		1,3-Cyclopentanedione, 2-acetyl...	140	C7H8O3	003859-39-0	5



Library Search Compound Report

Data File : C:\MSDCHEM\1\5973N\02FEB28\02716.D
 Acq Cn : 28 Feb 2002 2:40 pm
 Sample : 508 scan
 Misc : February 28, 2002
 MS Integration Params: LSCINT.P

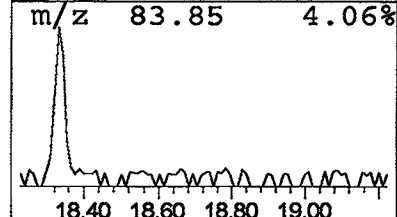
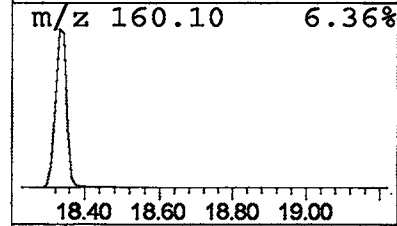
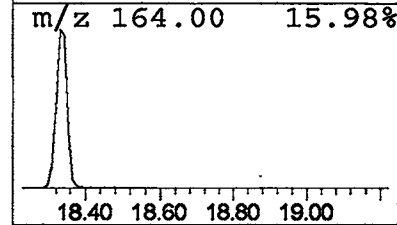
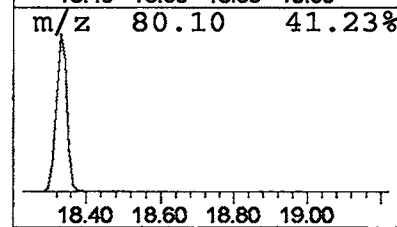
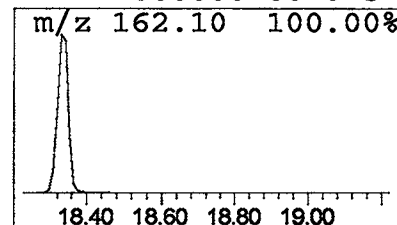
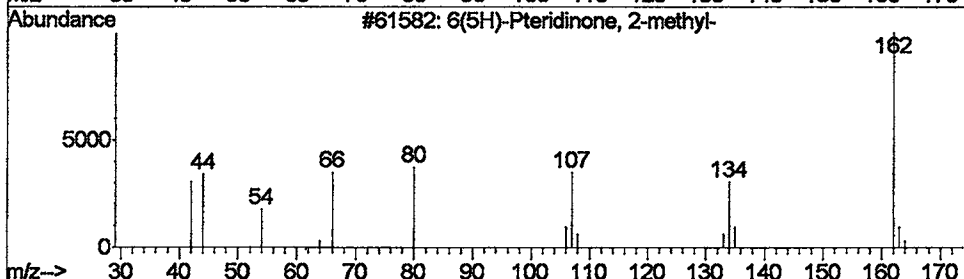
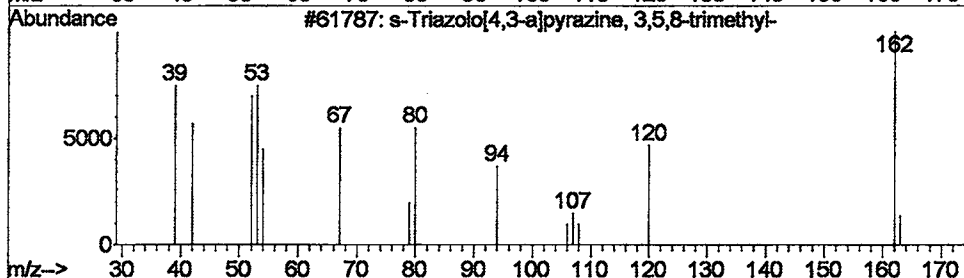
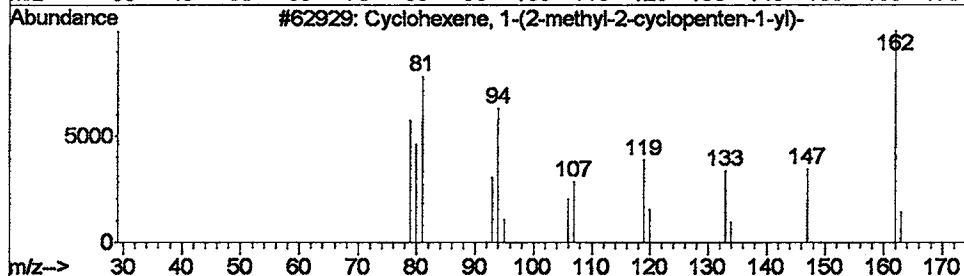
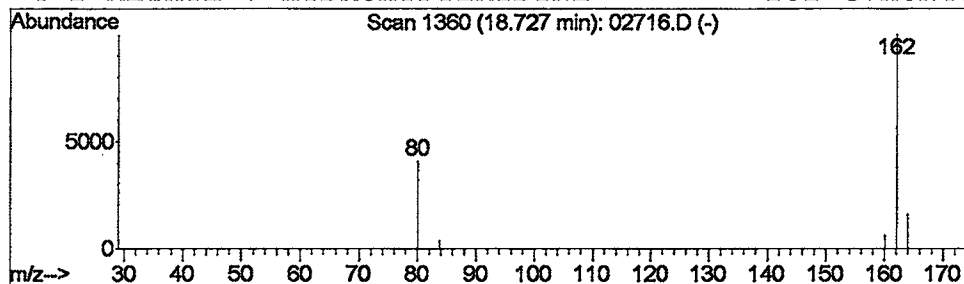
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Quant Method : C:\MSDCHEM\1...\HP022102.M (RTE Integrator)
 Title : 508 scan method
 Library : C:\DATABASE\WILEY7N.L

 Peak Number 9 Cyclohexene, 1-(2-methyl-2-... Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.73	0.03 ug/L	15611	Acenaphthene-d10	18.33

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclohexene, 1-(2-methyl-2-cyclo...	162	C12H18	058543-96-7	5
2			s-Triazolo[4,3-a]pyrazine, 3,5,8...	162	C8H10N4	019848-79-4	5
3			6(5H)-Pteridinone, 2-methyl-	162	C7H6N4O	016041-25-1	5
4			2-METHYL-7-HYDROXYPTERIDINE	162	C7H6N4O	000000-00-0	5



Library Search Compound Report

Data File : C:\MSDCHEM\1\5973N\02FEB28\02716.D
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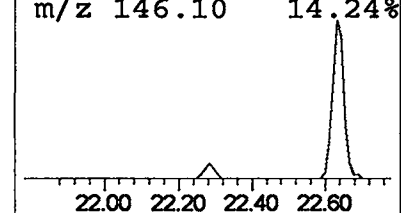
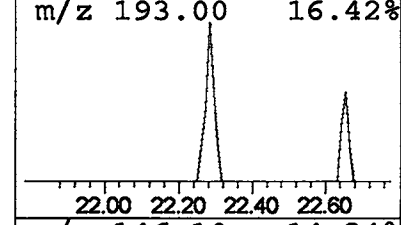
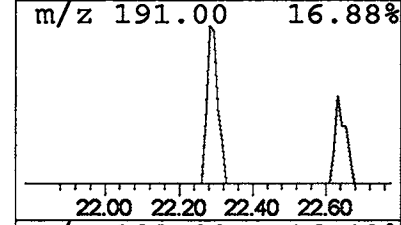
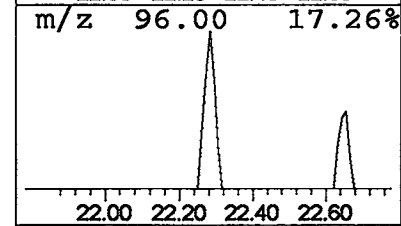
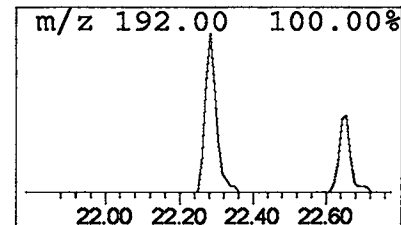
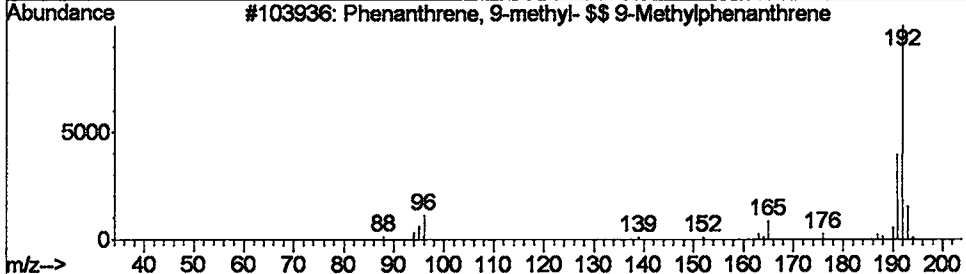
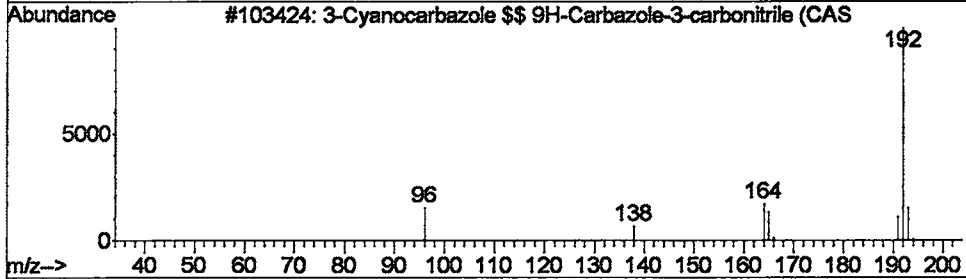
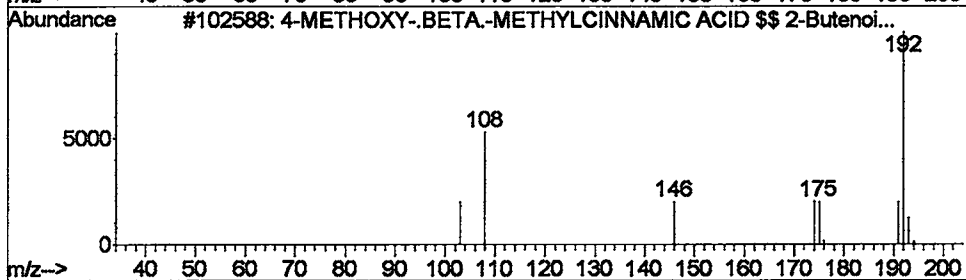
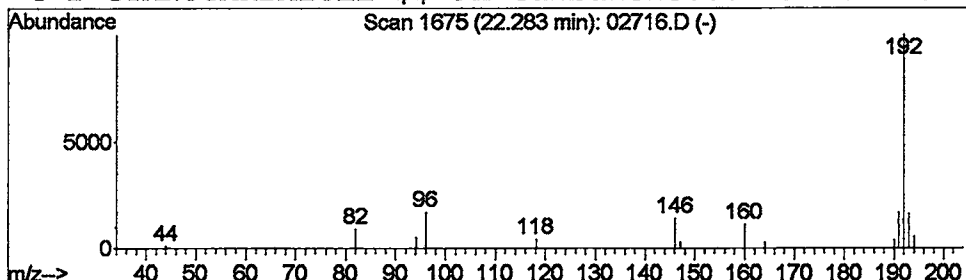
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Quant Method : C:\MSDCHEM\1...\HP022102.M (RTE Integrator)
 Title : 508 scan method
 Library : C:\DATABASE\WILEY7N.L

 Peak Number 10 4-METHOXY-.BETA.-METHYLCINN... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.28	0.12 ug/L	75603	Phenanthrene-d10	22.63

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	4-METHOXY-.BETA.-METHYLCINNAMIC ...	192	C11H12O3	019169-94-9	64
2		3-Cyanocarbazole \$\$ 9H-Carbazole...	192	C13H8N2	057102-93-9	59
3		Phenanthrene, 9-methyl- \$\$ 9-Met...	192	C15H12	000883-20-5	59
4		2-CYANOCARBAZOLE \$\$ 9H-Carbazole...	192	C13H8N2	057955-18-7	45



Library Search Compound Report

Data File : C:\MSDCHEM\1\5973N\02FEB28\02716.D
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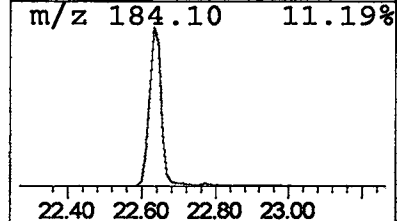
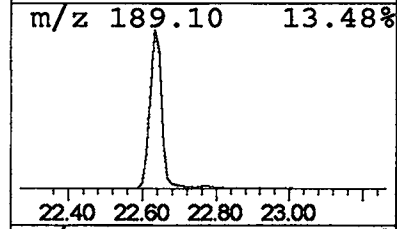
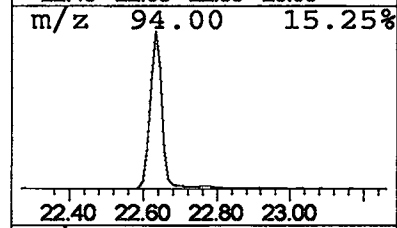
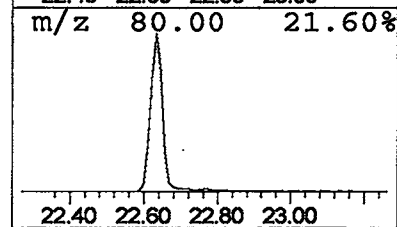
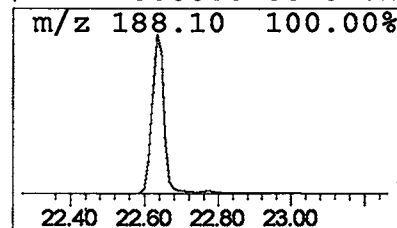
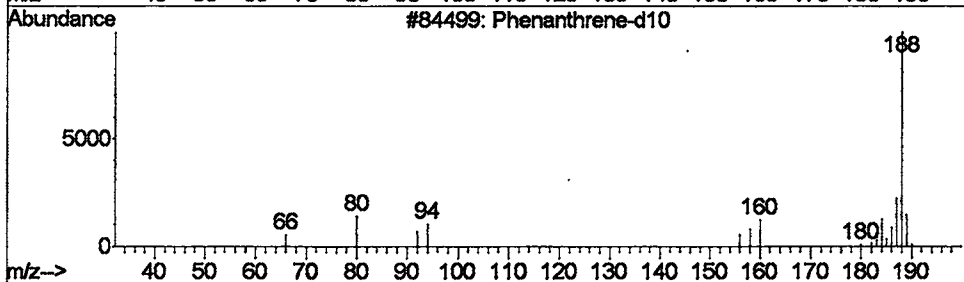
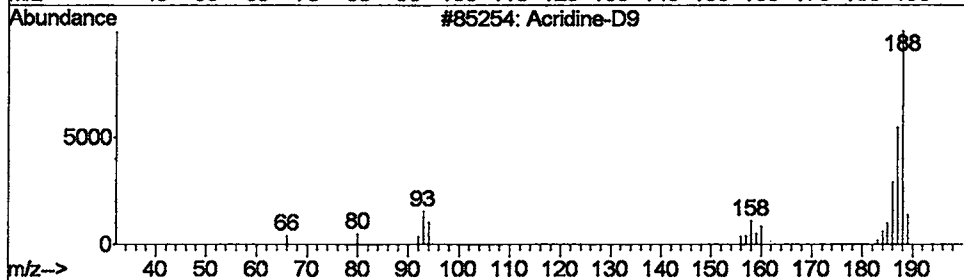
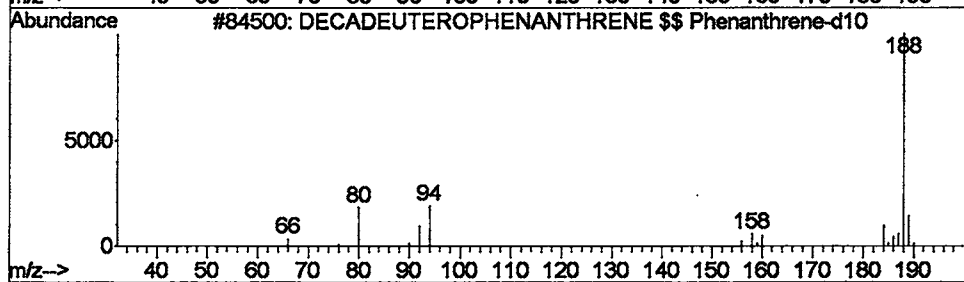
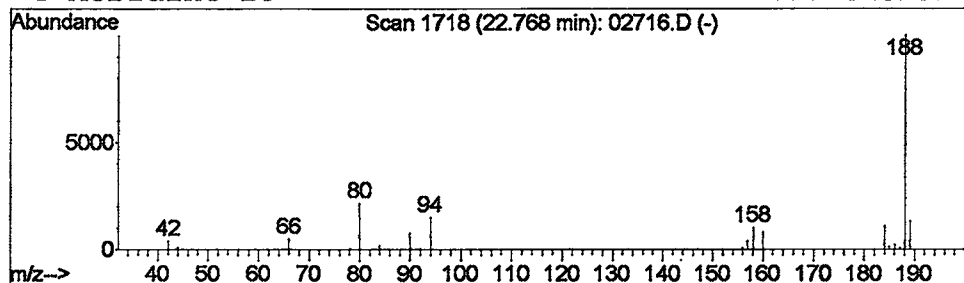
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Quant Method : C:\MSDCHEM\1...\HP022102.M (RTE Integrator)
 Title : 508 scan method
 Library : C:\DATABASE\WILEY7N.L

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.77	0.62 ug/L	396144	Phenanthrene-d10	22.63

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	DECADEUTEROPHENANTHRENE \$\$ Phena...	188	C14D10	001517-22-2	87
2		Acridine-D9	188	C13D9N	000000-00-0	80
3		Phenanthrene-d10	188	C14D10	001517-22-2	80
4		Acridine-D9	188	C13D9N	000000-00-0	72



Library Search Compound Report

Data File : C:\MSDCHEM\1\5973N\02FEB28\02716.D
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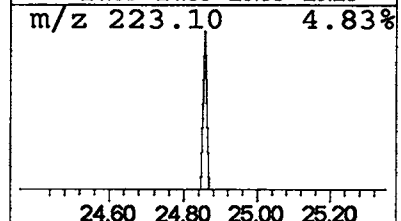
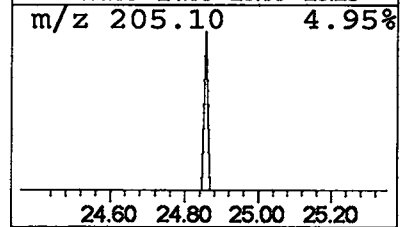
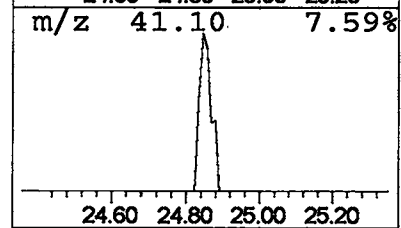
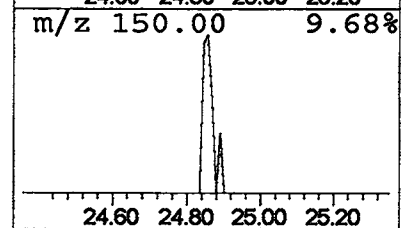
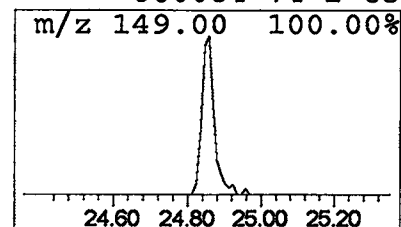
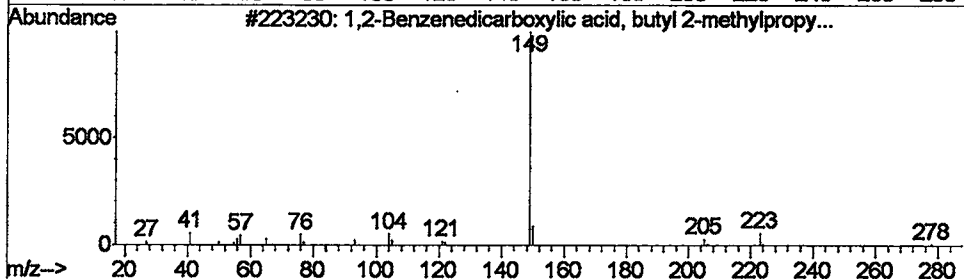
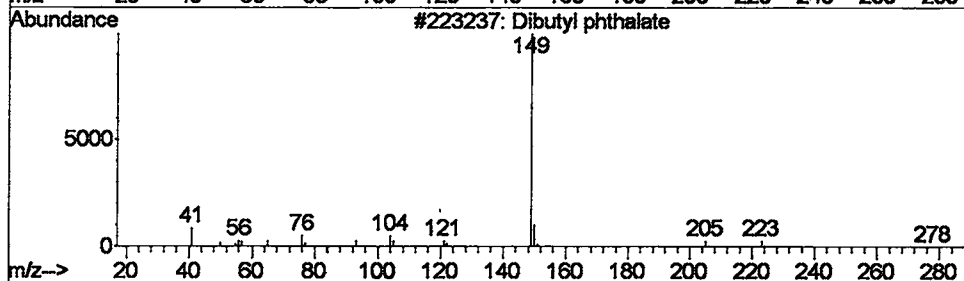
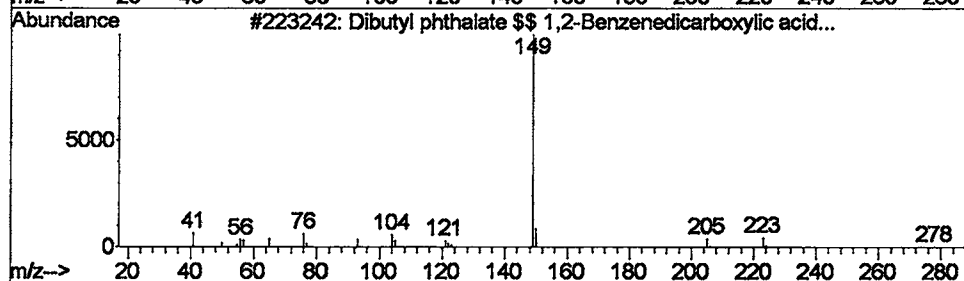
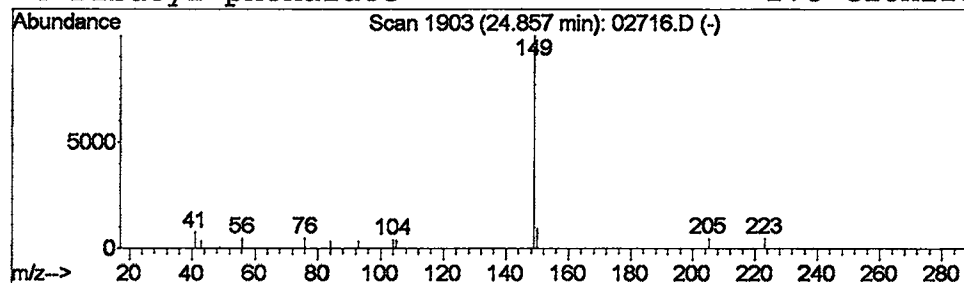
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 Inst : Instrumen
 Multiplr: 1.00

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

 Peak Number 12 Dibutyl phthalate \$\$ 1,2-Be... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.86	0.08 ug/L	47657	Phenanthrene-d10	22.63

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Dibutyl phthalate \$\$ 1,2-Benzene...	278	C16H22O4	000084-74-2	83
2			Dibutyl phthalate	278	C16H22O4	000084-74-2	83
3			1,2-Benzenedicarboxylic acid, bu...	278	C16H22O4	017851-53-5	83
4			Dibutyl phthalate	278	C16H22O4	000084-74-2	83



Library Search Compound Report

Data File : C:\MSDCHEM\1\5973N\02FEB28\02716.D
 Acq Cn : 28 Feb 2002 2:40 pm
 Sample : 508 scan
 Misc : February 28, 2002
 MS Integration Params: LSCINT.P

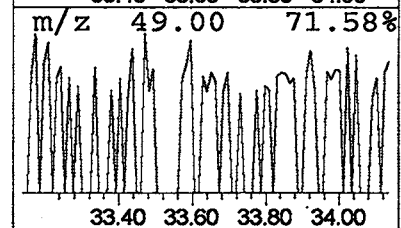
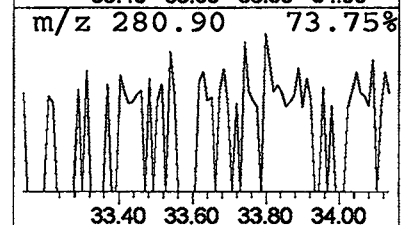
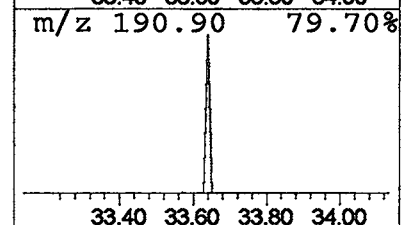
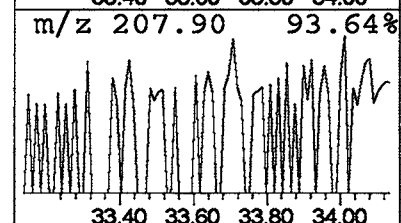
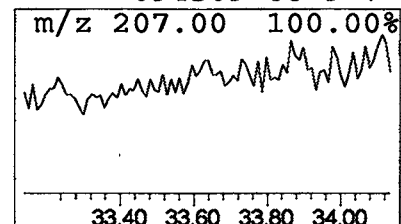
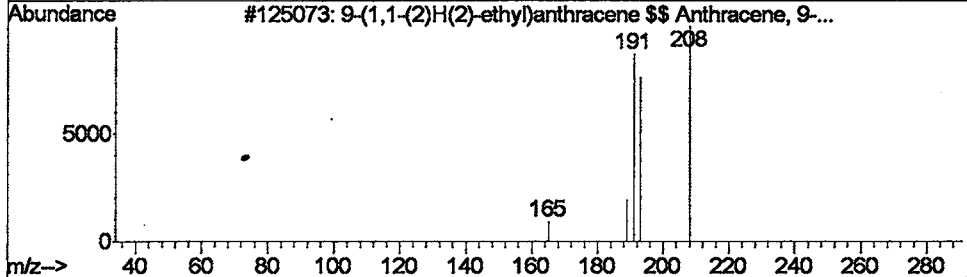
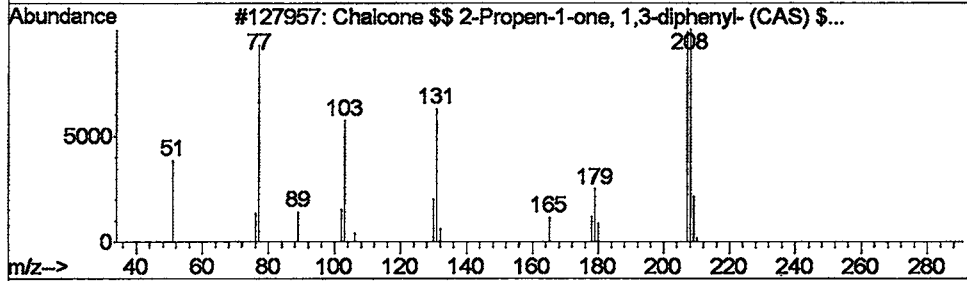
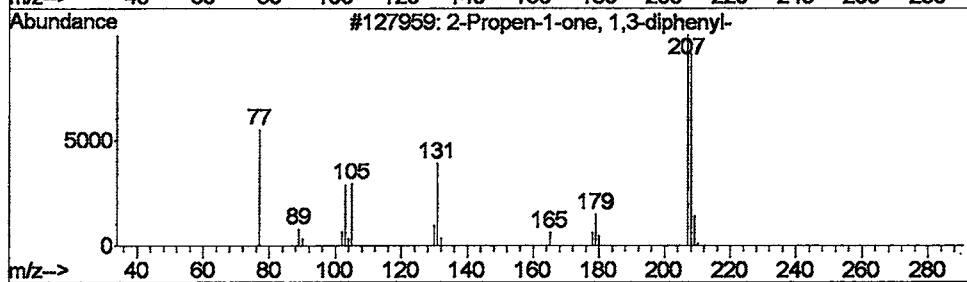
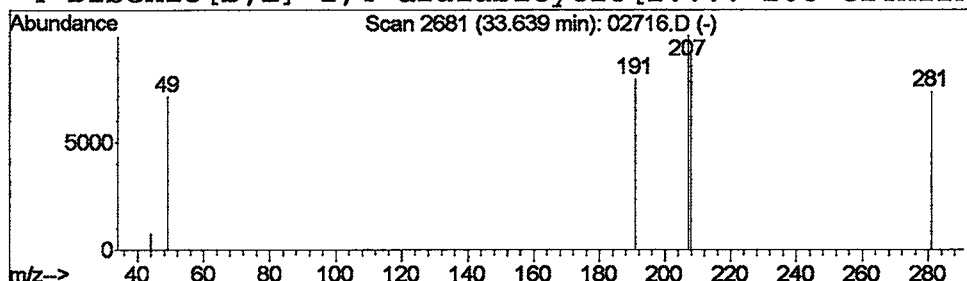
Vial: 1
 Operator:
 Inst : Instrumen
 Multiplr: 1.00

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

 Peak Number 13 2-Propen-1-one, 1,3-diphenyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
33.64	0.04 ug/L	17248	Perylene-d12	34.41

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Propen-1-one, 1,3-diphenyl-	208	C15H12O	000094-41-7	7
2			Chalcone \$\$ 2-Propen-1-one, 1,3-...	208	C15H12O	000094-41-7	7
3			9-(1,1-(2)H(2)-ethyl)anthracene ...	208	C16H12D2	122875-86-9	7
4			Dibenzo[b,E]-1,4-diazabicyclo[2....	208	C14H12N2	094509-68-9	7



Tentatively Identified Compound (LSC) summary

Operator ID: Date Acquired: 28 Feb 2002 2:40 pm
 Data File: C:\MSDCHEM\1\5973N\02FEB28\02716.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	0.1	ug/L	50459	1	9.72	7355020	20.0
ISOBUTYRALDEHYDE,...	3.93	0.1	ug/L	28562	1	9.72	7355020	20.0
2,2-dimethoxybutane	4.24	0.3	ug/L	128314	1	9.72	7355020	20.0
1-Propene, 3-meth...	5.02	0.1	ug/L	29102	1	9.72	7355020	20.0
Imidazole-2,4,5-d...	5.72	0.0	ug/L	13724	1	9.72	7355020	20.0
Acetic acid, 3-[1...	5.95	0.3	ug/L	107145	1	9.72	7355020	20.0
2-Cyclopenten-1-o...	13.38	0.0	ug/L	25450	2	13.20	10781000	20.0
3-Hydroxy-2,2,4-t...	16.55	0.0	ug/L	21076	3	18.33	11831500	20.0
Cyclohexene, 1-(2...	18.73	0.0	ug/L	15611	3	18.33	11831500	20.0
4-METHOXY-.BETA.-...	22.28	0.1	ug/L	75603	4	22.63	12694200	20.0
DECADEUTEROPHENAN...	22.77	0.6	ug/L	396144	4	22.63	12694200	20.0
Dibutyl phthalate...	24.86	0.1	ug/L	47657	4	22.63	12694200	20.0
2-Propen-1-one, 1...	33.64	0.0	ug/L	17248	6	34.41	8473810	20.0