

User: Vinci, David L
Westchester Dept. of Labs & Research
Date: 02/26/2002 Time: 11:10:02

Sample: AE00842
Analysis: \$GCMS GCMS Scan For Unknowns
Units: ug
Started: 2/26/02 11:09 Ended: 2/26/02 11:09 Analyst: DLV
Page: 1

	Component Name	Viol	Result	Qualifier	MDL
1	Other unknown compounds		.		
2	Surr_2,4-DDD		114		5.0

Comments for Sample AE00842 - Analysis \$GCMS

Westchester Dept. of Labs & Research
User: Vinci, David L
Date: 02-26-2002 Time: 11:11:01

The GCMS scan, from 35 to 550 amu, does not reveal the presence of unusual compounds. The recoveries for 21 of the 43 compounds in the LCS Standard were above their acceptance ranges.

Data File : C:\MSDCHEM\1\5973N\02FEB07\00842.D

Vial: 6

Acq On : 7 Feb 2002 6:34 pm

Operator:

Sample : 508 scan

Inst : Instrumen

Misc : February 7, 2002

Multiplr: 1.00

MS Integration Params: SPLIT.P

Quant Time: Feb 20 15:39:46 2002

Quant Results File: HP020702.RES

Quant Method : C:\MSDCHEM\1...\HP020702.M (RTE Integrator)

Title : 508 scan method

Last Update : Wed Feb 20 11:48:57 2002

Response via : Initial Calibration

DataAcq Meth : SCAN508

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev (Min)
1) 1,4-Dichlorobenzene-d4	9.71	152	1443116	20.00	ug/L	0.00
10) Naphthalene-d8	13.19	136	6023418	20.00	ug/L	-0.01
17) Acenaphthene-d10	18.34	164	3108615	20.00	ug/L	0.00
29) Phenanthrene-d10	22.63	188	5071337	20.00	ug/L	-0.01
38) Chrysene-d12	30.49	240	4438090	20.00	ug/L	-0.03
44) Perylene-d12	34.42	264	3137991	20.00	ug/L	-0.03

System Monitoring Compounds

37) 2,4-DDD	27.73	235	476239	5.70	ug/L	0.00
Spiked Amount	5.000		Recovery	=	114.00%	

Target Compounds

						Qvalue
20) Dimethylphthalate	18.34	163	489420	2.60	ug/L	# 58
21) 2,6-Dinitrotoluene	18.34	165	392016	9.00	ug/L	# 27

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

00842.D HP020702.M Wed Feb 20 15:39:47 2002

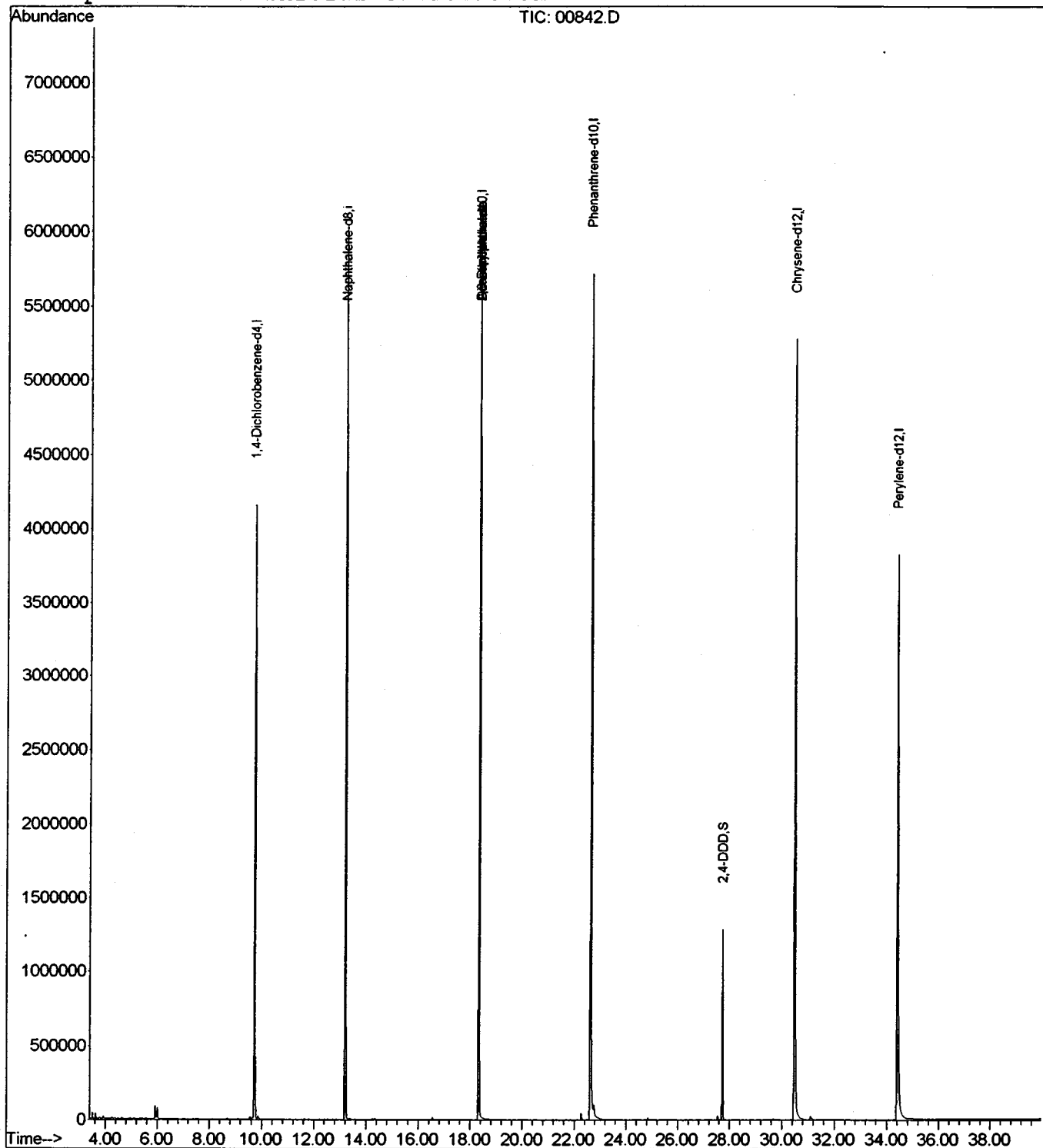
Page 1

Data File : C:\MSDCHEM\1\5973N\02FEB07\00842.D
 Acq On : 7 Feb 2002 6:34 pm
 Sample : 508 scan
 Misc : February 7, 2002
 MS Integration Params: SPLIT.P
 Quant Time: Feb 20 15:39 2002

Vial: 6
 Operator:
 Inst : Instrumen
 Multiplr: 1.00

Quant Results File: HP020702.R

Method : C:\MSDCHEM\1\METHODS\METHODS\HP020702.M (RTE Integrator)
 Title : 508 scan method
 Last Update : Wed Feb 20 11:48:57 2002
 Response via : Initial Calibration



Data File : C:\MSDCHEM\1\5973N\02FEB07\00842.D
 Acq On : 7 Feb 2002 6:34 pm
 Sample : 508 scan
 Misc : February 7, 2002
 MS Integration Params: LSCINT.P

Vial: 6
 Operator:
 Inst : Instrumen
 Multiplr: 1.00

Method : C:\MSDCHEM\1\METHODS\METHODS\HP020702.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

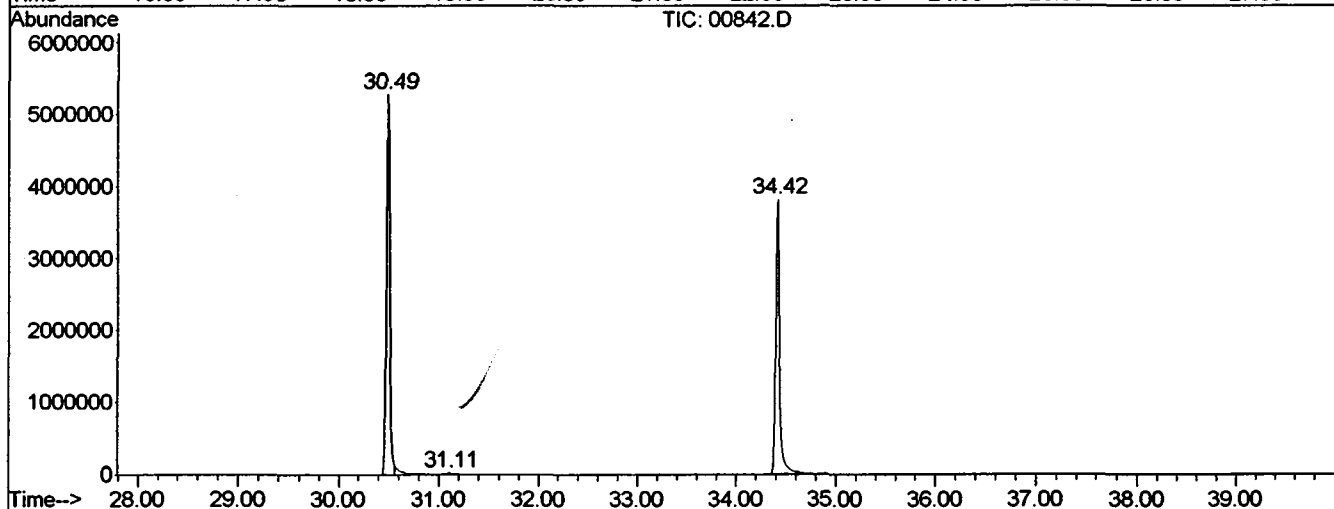
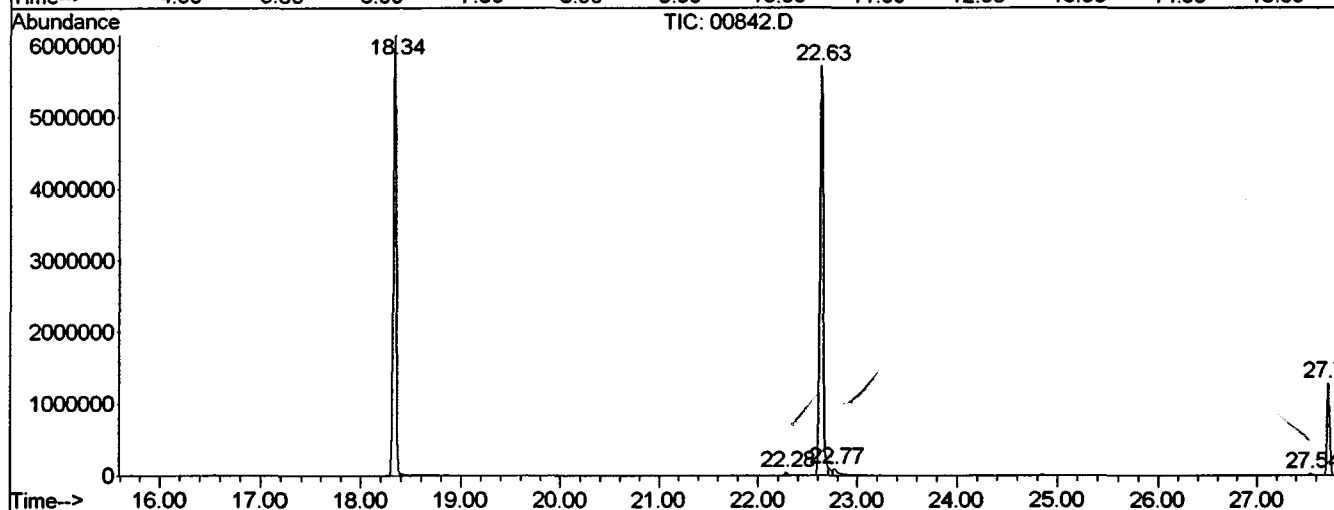
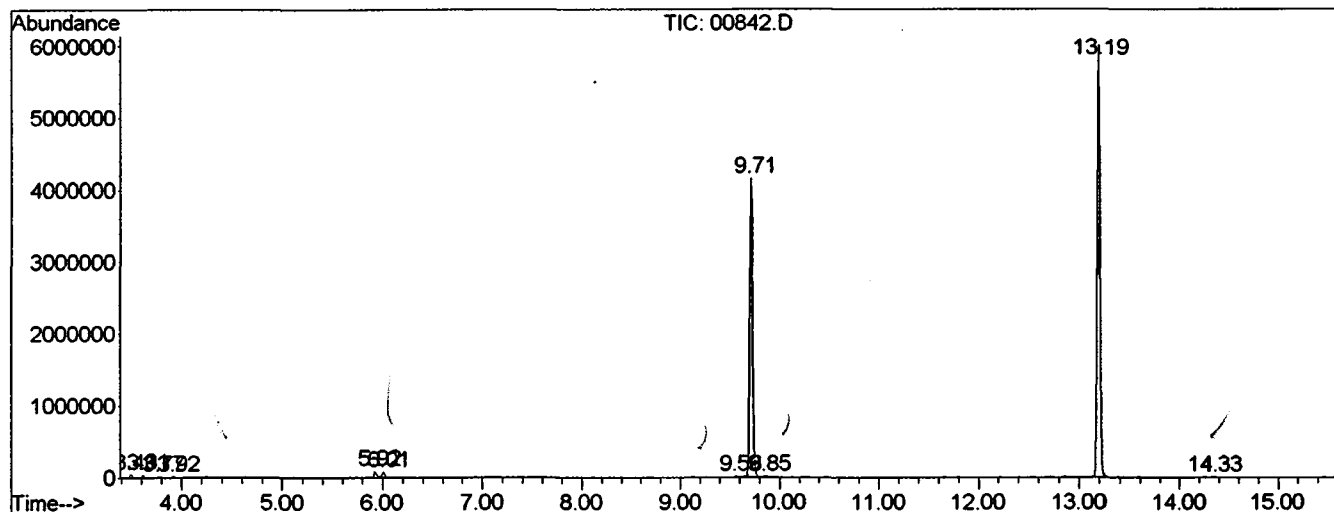
If leading or trailing edge < 100 prefer < Baseline drop else tangent >
 Peak separation: 5

Signal : TIC

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.488	5	10	15	rBV	45072	79931	0.60%	0.111%
2	3.613	19	21	25	rBV	40695	58894	0.44%	0.082%
3	3.773	31	35	45	rVV3	9795	39480	0.30%	0.055%
4	3.922	45	48	53	rVB2	17409	32960	0.25%	0.046%
5	5.920	219	223	227	rBV	92010	172569	1.30%	0.240%
6	6.011	227	231	235	rVB	75111	147393	1.11%	0.205%
7	9.562	537	542	547	rBV	17727	44022	0.33%	0.061%
8	9.710	551	555	561	rVV	4155906	8159577	61.26%	11.327%
9	9.847	565	567	575	rBV	20182	52151	0.39%	0.072%
10	13.192	855	860	865	rBV	6001052	11412355	85.68%	15.843%
11	14.334	957	960	969	rVB4	8915	31120	0.23%	0.043%
12	18.341	1305	1311	1315	rBV	6147701	12308829	92.41%	17.088%
13	22.280	1653	1656	1661	rVV	44707	89734	0.67%	0.125%
14	22.634	1681	1687	1693	rBV2	5715375	13319274	100.00%	18.490%
15	22.771	1697	1699	1729	rVB	100945	451044	3.39%	0.626%
16	27.543	2111	2117	2127	rBV	28590	67550	0.51%	0.094%
17	27.726	2127	2133	2139	rBV	1288903	2311270	17.35%	3.209%
18	30.489	2369	2375	2381	rBV2	5278449	12646256	94.95%	17.556%
19	31.105	2419	2429	2447	rVB5	26173	118534	0.89%	0.165%
20	34.416	2713	2719	2753	rBV	3821865	10490437	78.76%	14.563%

Sum of corrected areas: 72033380

File : C:\MSDCHEM\1\5973N\02FEB07\00842.D
 Operator :
 Acquired : 7 Feb 2002 6:34 pm using AcqMethod SCAN508
 Instrument : Instrumen
 Sample Name: 508 scan
 Misc Info : February 7, 2002
 Vial Number: 6
 Quant File :HP020702.RES (RTE Integrator)



Data File : C:\MSDCHEM\1\5973N\02FEB07\00842.D
 Acq On : 7 Feb 2002 6:34 pm
 Sample : 508 scan
 Misc : February 7, 2002
 MS Integration Params: LSCINT.P

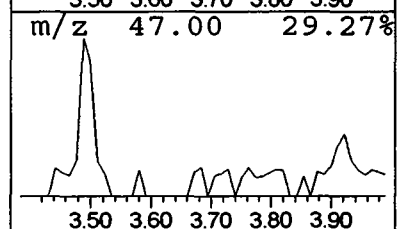
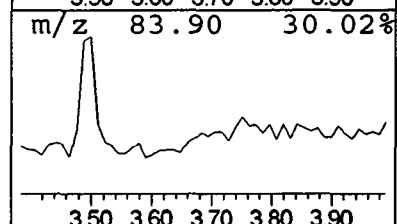
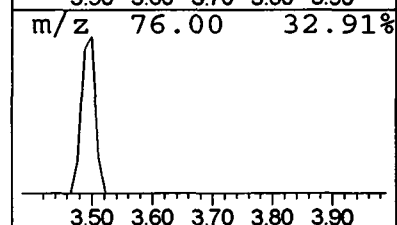
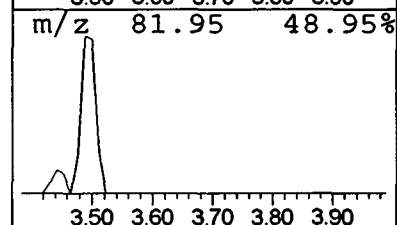
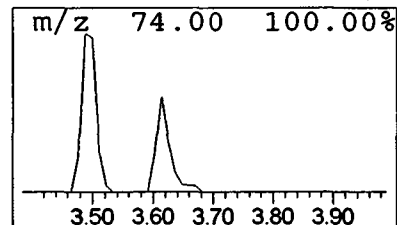
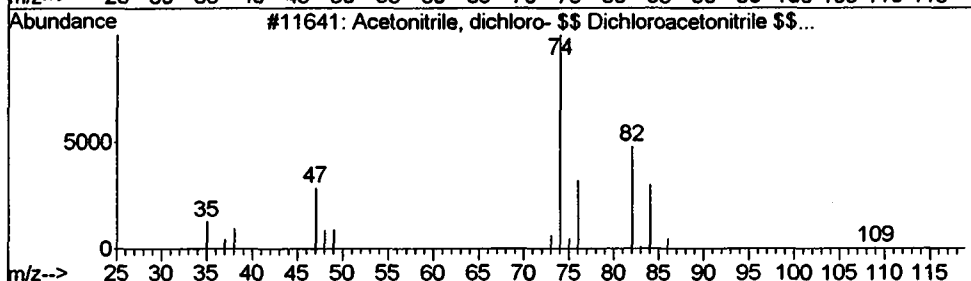
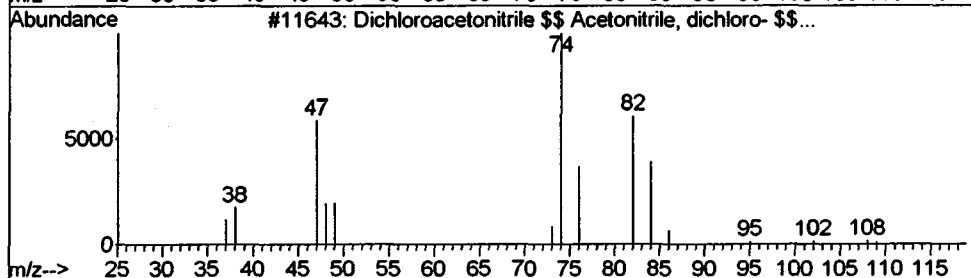
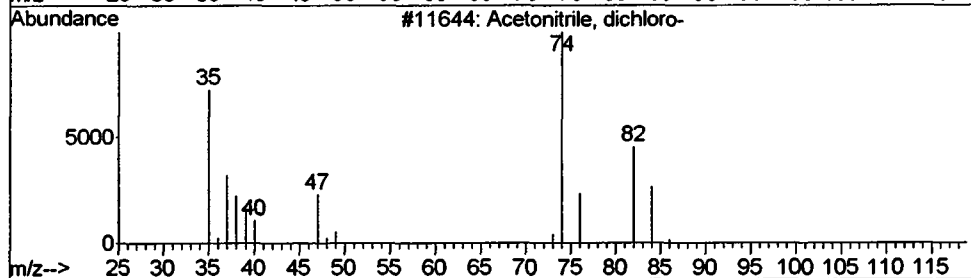
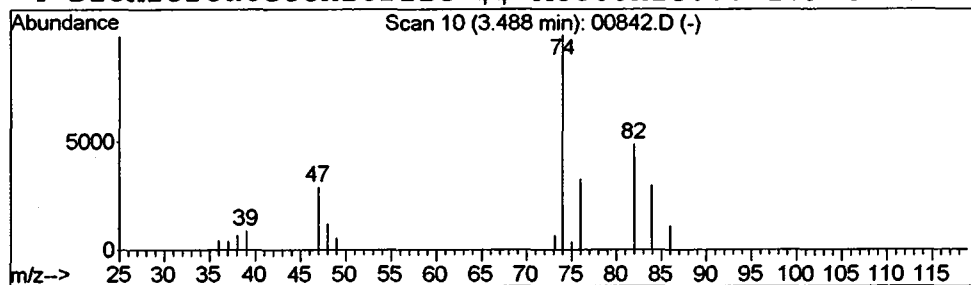
Vial: 6
 Operator:
 Inst : Instrumen
 Multiplr: 1.00

Quant Method : C:\MSDCHEM\1...\HP020702.M (RTE Integrator)
 Title : 508 scan method
 Library : C:\DATABASE\WILEY7N.L

 Peak Number 1 Acetonitrile, dichloro- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.49	0.20 ug/L	79931	1,4-Dichlorobenzene-d4	9.71

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Acetonitrile, dichloro-	109	C2HCl2N	003018-12-0	78
2		Dichloroacetonitrile \$\$ Acetonit...	109	C2HCl2N	003018-12-0	74
3		Acetonitrile, dichloro- \$\$ Dichl...	109	C2HCl2N	003018-12-0	64
4		Dichloroacetonitrile \$\$ Acetonit...	109	C2HCl2N	003018-12-0	53



Data File : C:\MSDCHEM\1\5973N\02FEB07\00842.D
 Acq On : 7 Feb 2002 6:34 pm
 Sample : 508 scan
 Misc : February 7, 2002
 MS Integration Params: LSCINT.P

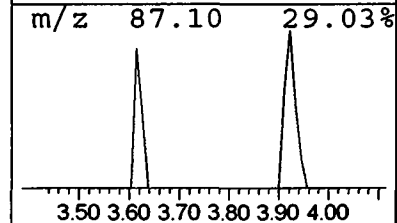
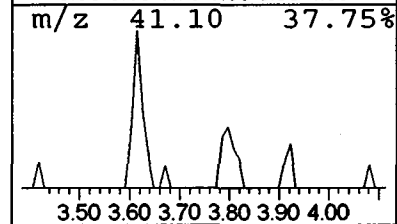
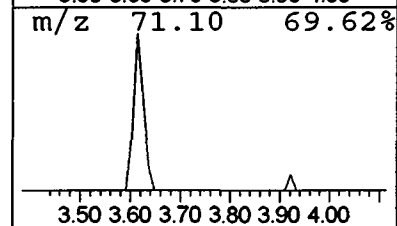
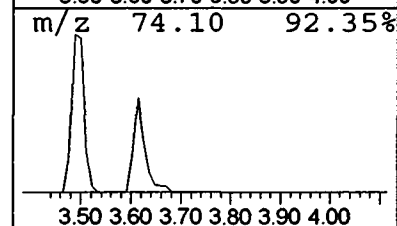
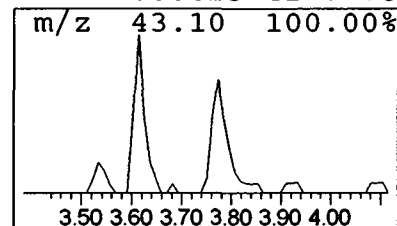
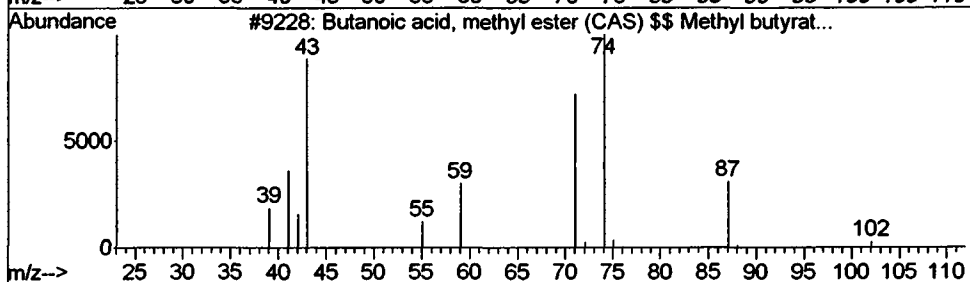
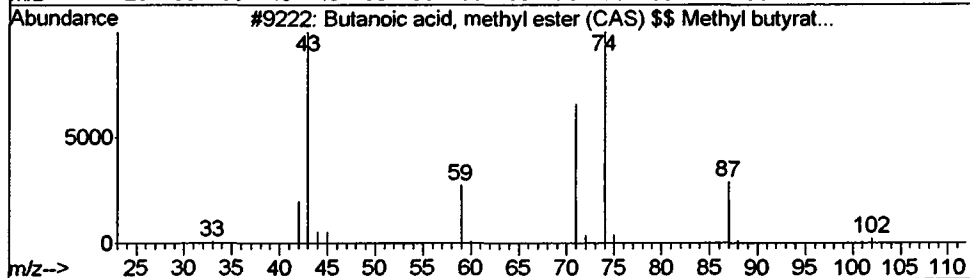
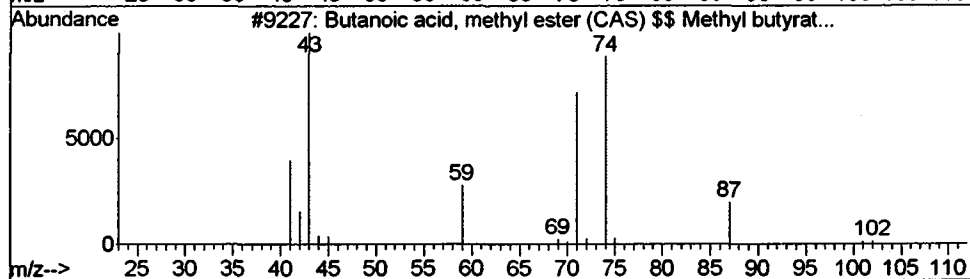
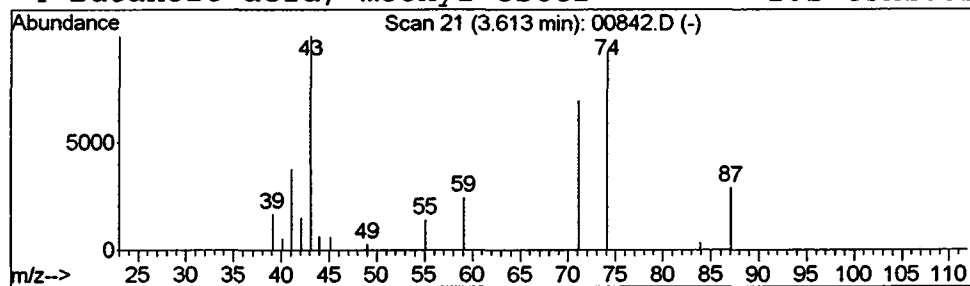
Vial: 6
 Operator:
 Inst : Instrumen
 Multiplr: 1.00

Quant Method : C:\MSDCHEM\1...\HP020702.M (RTE Integrator)
 Title : 508 scan method
 Library : C:\DATABASE\WILEY7N.L

 Peak Number 2 Butanoic acid, methyl ester... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.61	0.14 ug/L	58894	1,4-Dichlorobenzene-d4	9.71

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	78
4			Butanoic acid, methyl ester	102	C5H10O2	000623-42-7	78



Data File : C:\MSDCHEM\1\5973N\02FEB07\00842.D
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 Sample : 508 scan
 Misc : February 7, 2002
 MS Integration Params: LSCINT.P

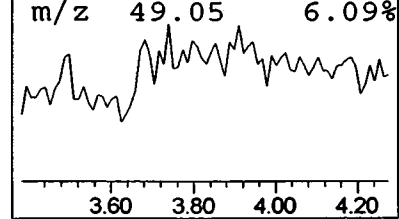
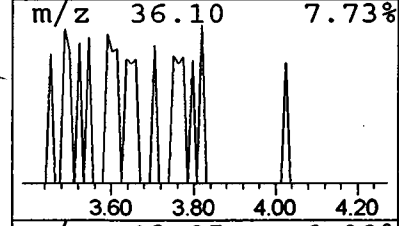
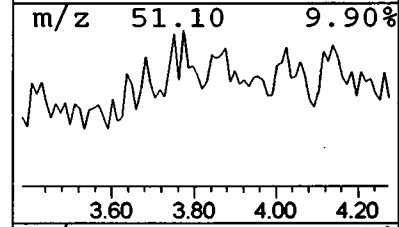
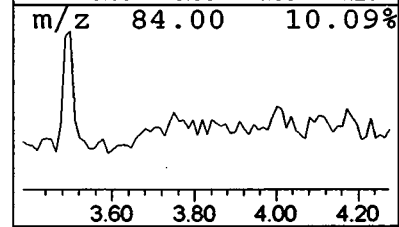
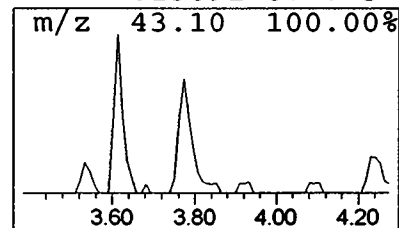
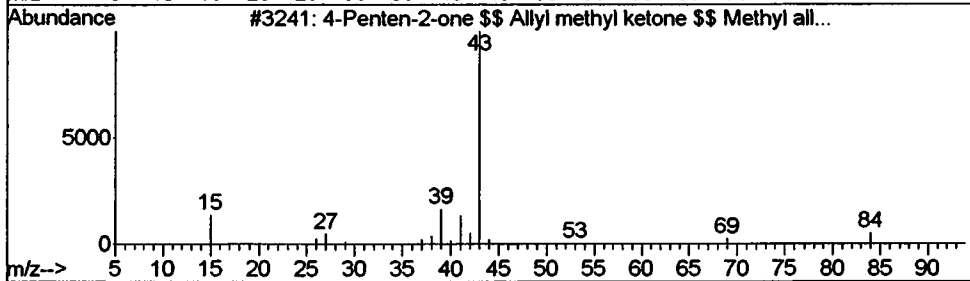
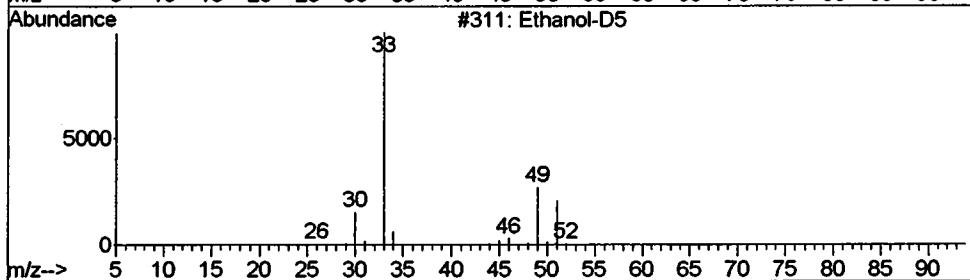
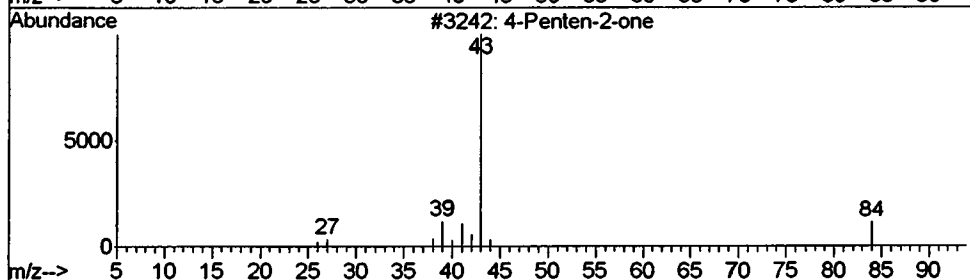
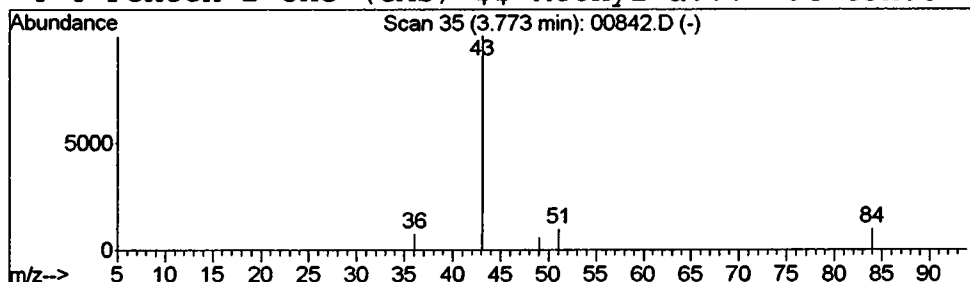
Vial: 6
 Operator:
 Inst : Instrumen
 Multiplr: 1.00

Quant Method : C:\MSDCHEM\1...\HP020702.M (RTE Integrator)
 Title : 508 scan method
 Library : C:\DATABASE\WILEY7N.L

 Peak Number 3 4-Penten-2-one Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.77	0.10 ug/L	39480	1,4-Dichlorobenzene-d4	9.71

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	4-Penten-2-one	84	C5H8O	013891-87-7	3
2	Ethanol-D5	51	C2HD5O	000000-00-0	3
3	4-Penten-2-one \$\$ Allyl methyl k...	84	C5H8O	013891-87-7	3
4	4-Penten-2-one (CAS) \$\$ Methyl a...	84	C5H8O	013891-87-7	3



Data File : C:\MSDCHEM\1\5973N\02FEB07\00842.D
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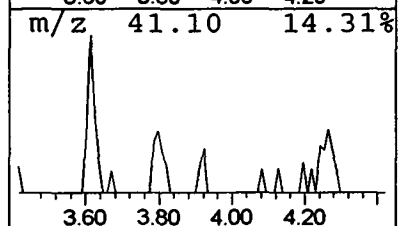
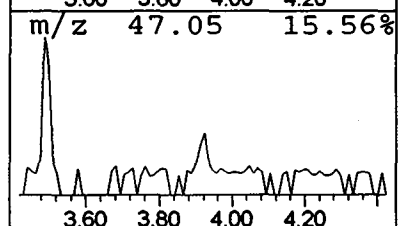
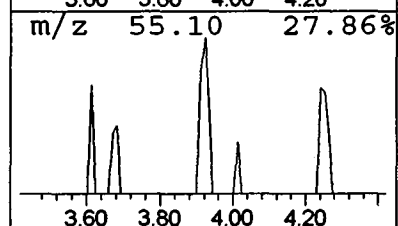
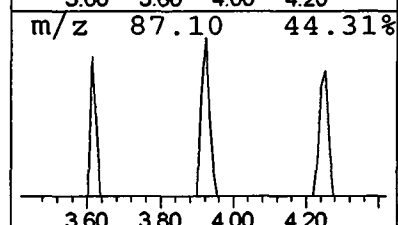
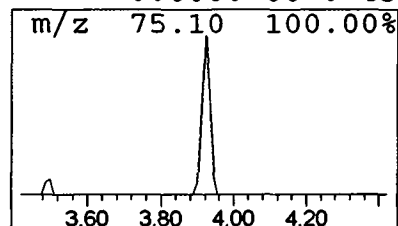
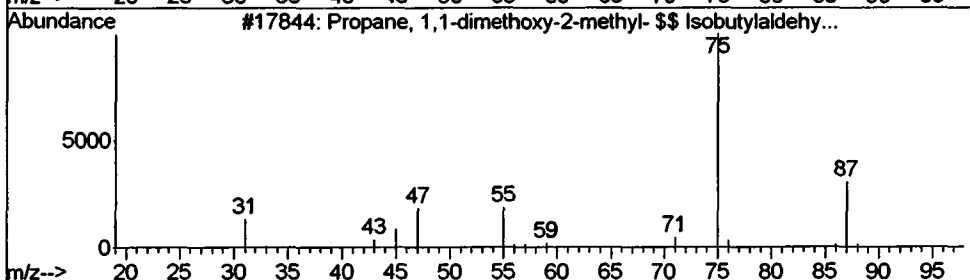
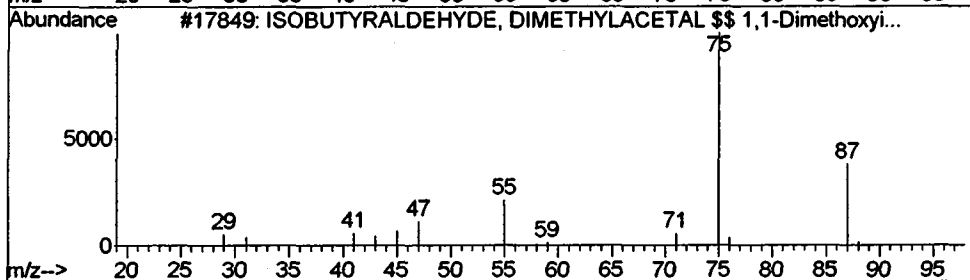
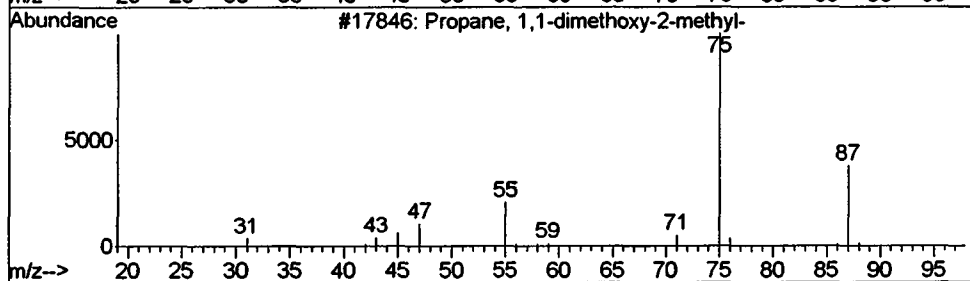
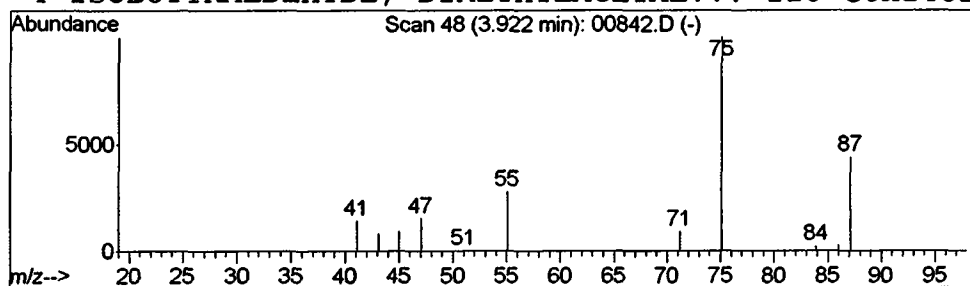
Vial: 6
 Operator:
 Inst : Instrumen
 Multiplr: 1.00

Quant Method : C:\MSDCHEM\1...\HP020702.M (RTE Integrator)
 Title : 508 scan method
 Library : C:\DATABASE\WILEY7N.L

 Peak Number 4 Propane, 1,1-dimethoxy-2-me... Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.92	0.08 ug/L	32960	1,4-Dichlorobenzene-d4	9.71

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Propane, 1,1-dimethoxy-2-methyl-	118	C6H14O2	041632-89-7	83
2		ISOBUTYRALDEHYDE, DIMETHYLACETAL...	118	C6H14O2	000000-00-0	83
3		Propane, 1,1-dimethoxy-2-methyl-	118	C6H14O2	041632-89-7	43
4		ISOBUTYRALDEHYDE, DIMETHYLACETAL...	118	C6H14O2	000000-00-0	43



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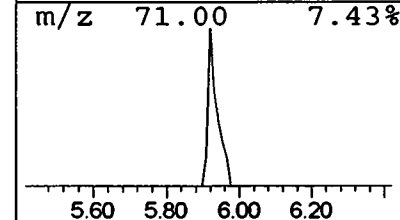
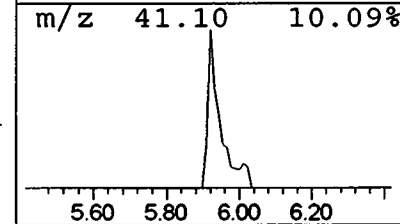
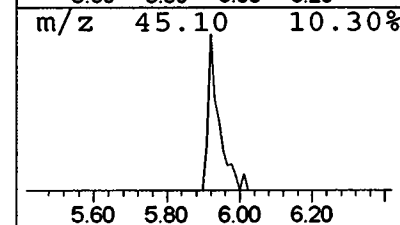
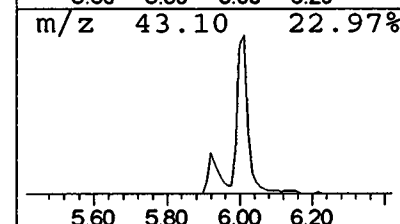
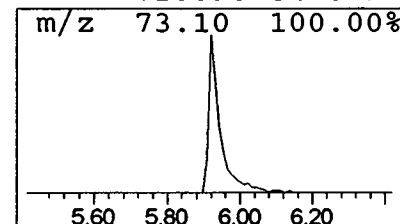
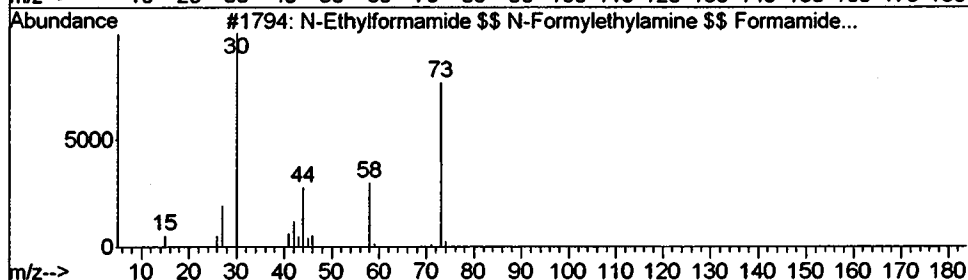
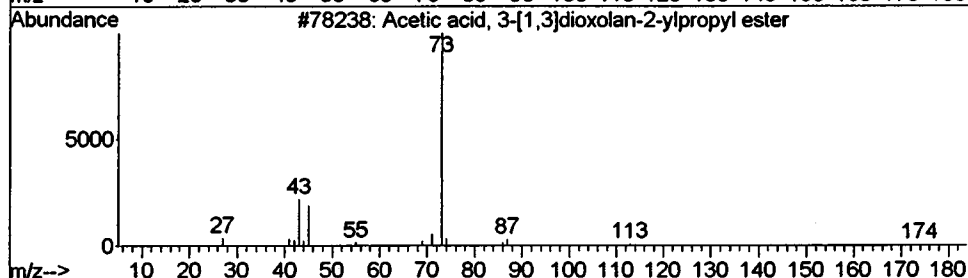
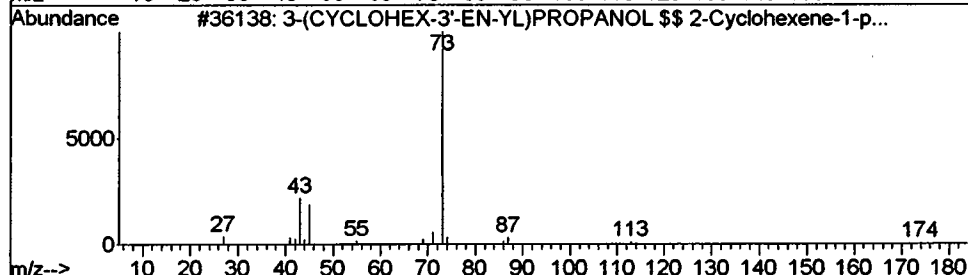
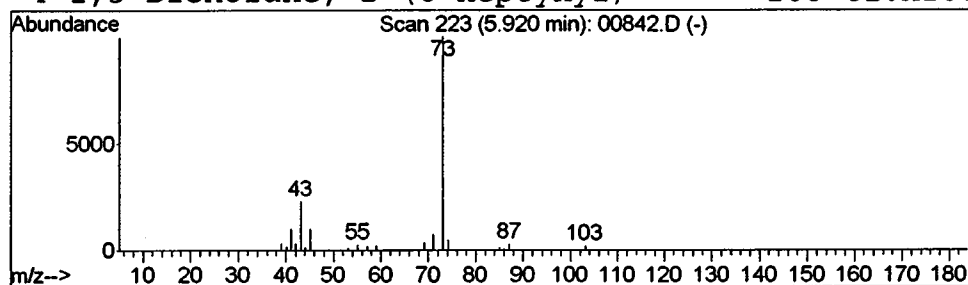
Vial: 6
 Operator:
 Inst : Instrumen
 Multiplr: 1.00

Quant Method : C:\MSDCHEM\1...\HP020702.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.
5.92	0.42 ug/L	172569	1,4-Dichlorobenzene-d4	9.71

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			3-(CYCLOHEX-3'-EN-YL)PROPANOL \$\$...	174	C8H14O4	015745-87-6	40
2			Acetic acid, 3-[1,3]dioxolan-2-y...	174	C8H14O4	000000-00-0	40
3			N-Ethylformamide \$\$ N-Formylethy...	73	C3H7NO	000627-45-2	9
4			1,3-Dioxolane, 2-(6-heptynyl)-	168	C10H16O2	025595-82-8	9



Data File : C:\MSDCHEM\1\5973N\02FEB07\00842.D
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 Misc : February 7, 2002
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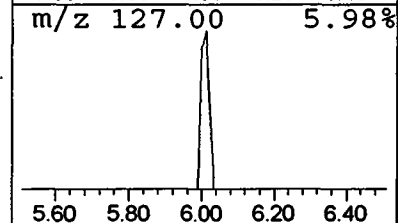
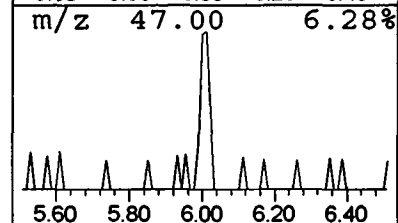
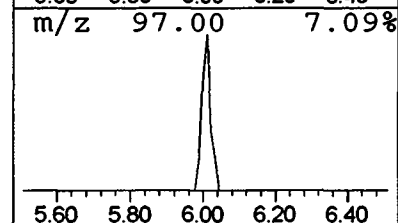
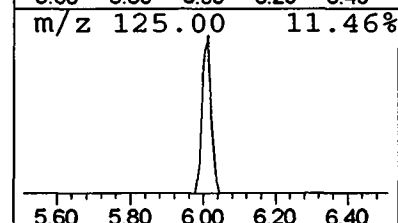
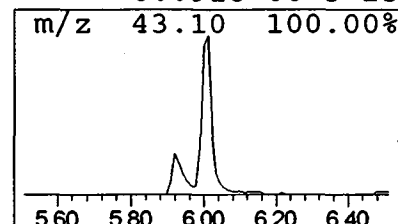
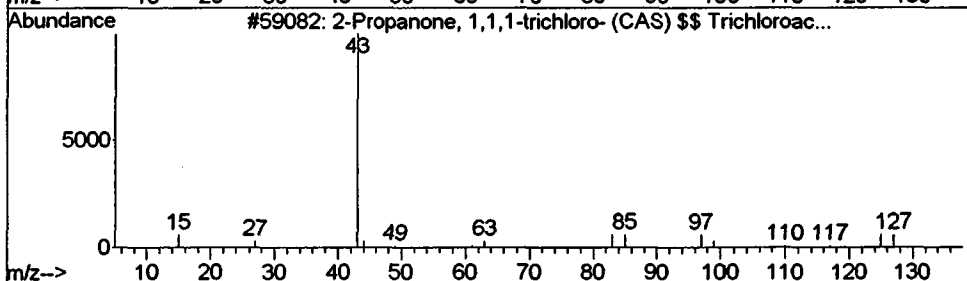
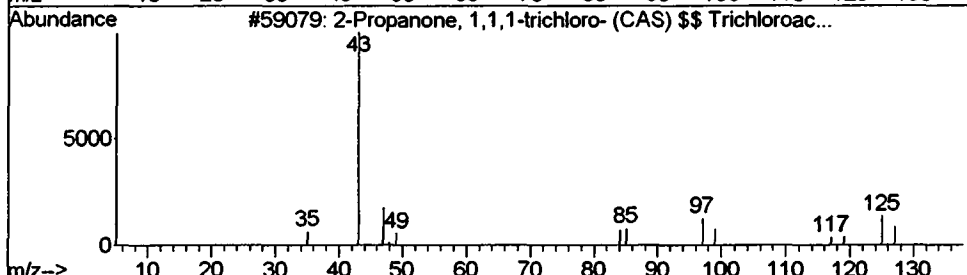
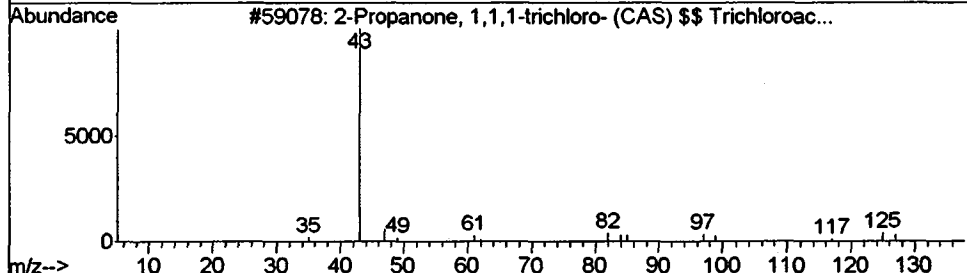
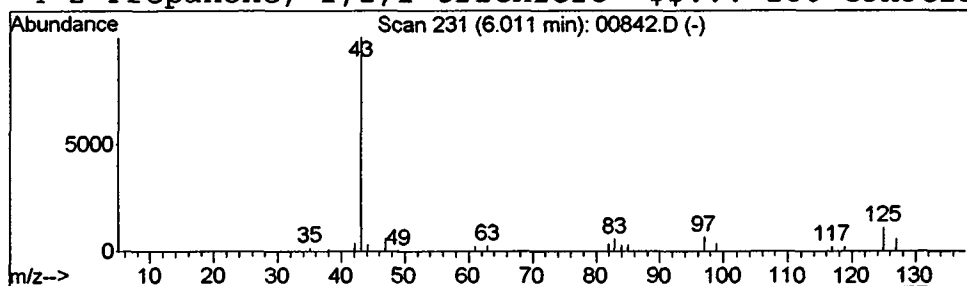
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 Inst : Instrumen
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Quant Method : C:\MSDCHEM\1...\HP020702.M (RTE Integrator)
 Title : 508 scan method
 Library : C:\DATABASE\WILEY7N.L

 Peak Number 6 2-Propanone, 1,1,1-trichlor... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.01	0.36 ug/L	147393	1,4-Dichlorobenzene-d4	9.71

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Propanone, 1,1,1-trichloro- (C...	160	C3H3Cl3O	000918-00-3	74
2			2-Propanone, 1,1,1-trichloro- (C...	160	C3H3Cl3O	000918-00-3	45
3			2-Propanone, 1,1,1-trichloro- (C...	160	C3H3Cl3O	000918-00-3	45
4			2-Propanone, 1,1,1-trichloro- \$\$...	160	C3H3Cl3O	000918-00-3	25



Data File : C:\MSDCHEM\1\5973N\02FEB07\00842.D
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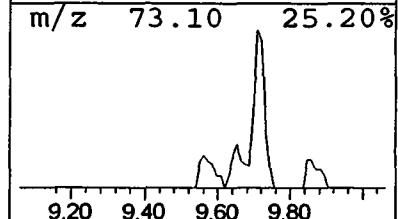
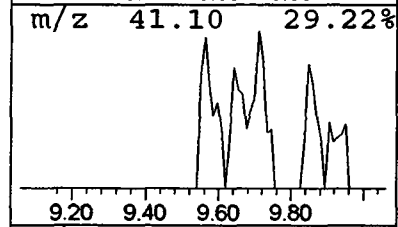
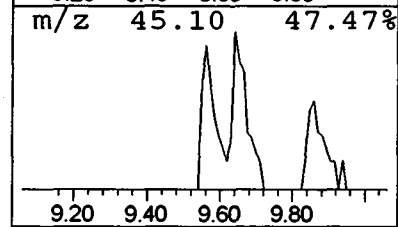
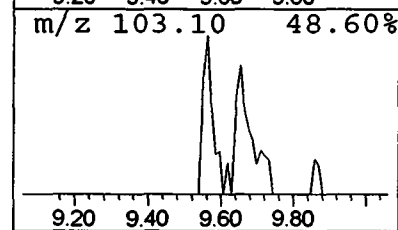
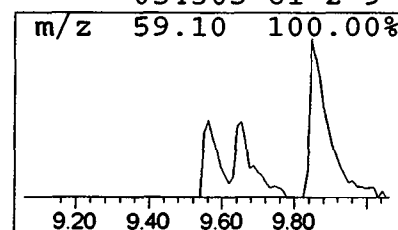
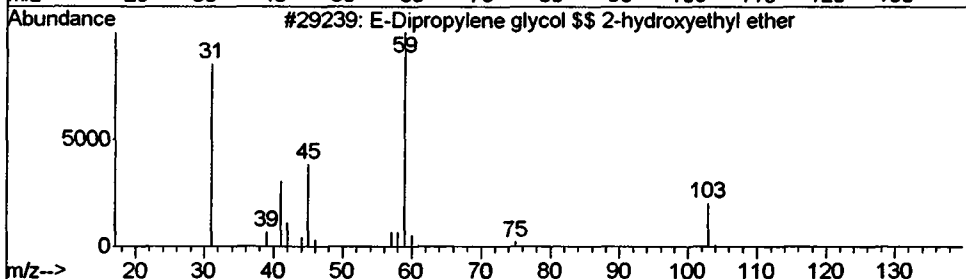
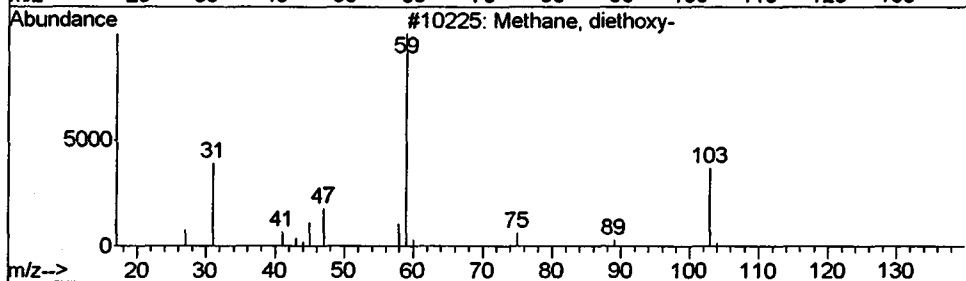
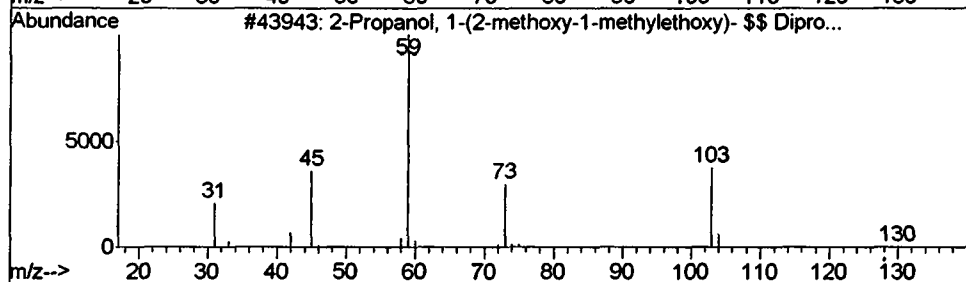
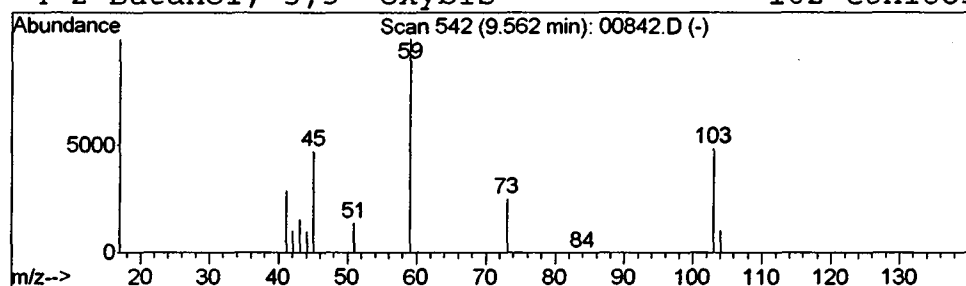
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 Multiplr: 1.00

Quant Method : C:\MSDCHEM\1...\HP020702.M (RTE Integrator)
 Title : 508 scan method
 Library : C:\DATABASE\WILEY7N.L

 Peak Number 7 2-Propanol, 1-(2-methoxy-1-... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.56	0.11 ug/L	44022	1,4-Dichlorobenzene-d4	9.71

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Propanol, 1-(2-methoxy-1-methy...	148	C7H16O3	020324-32-7	39
2	Methane, diethoxy-	104	C5H12O2	000462-95-3	9
3	E-Dipropylene glycol \$\$ 2-hydrox...	134	C6H14O3	000000-00-0	9
4	2-Butanol, 3,3'-oxybis-	162	C8H18O3	054305-61-2	9



Data File : C:\MSDCHEM\1\5973N\02FEB07\00842.D
 Acq On : 7 Feb 2002 6:34 pm
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 Misc : February 7, 2002
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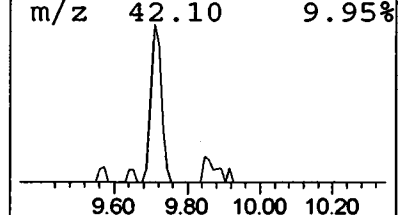
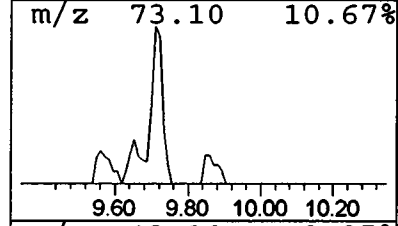
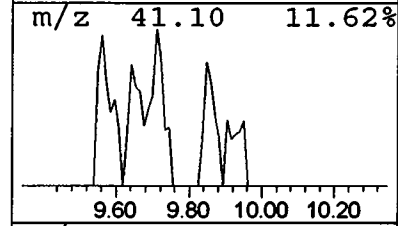
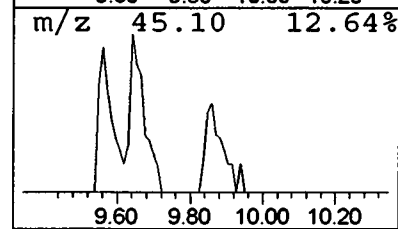
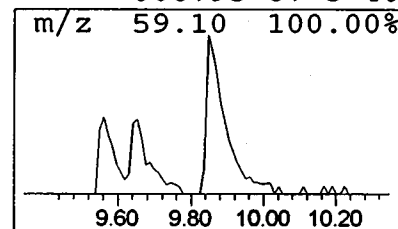
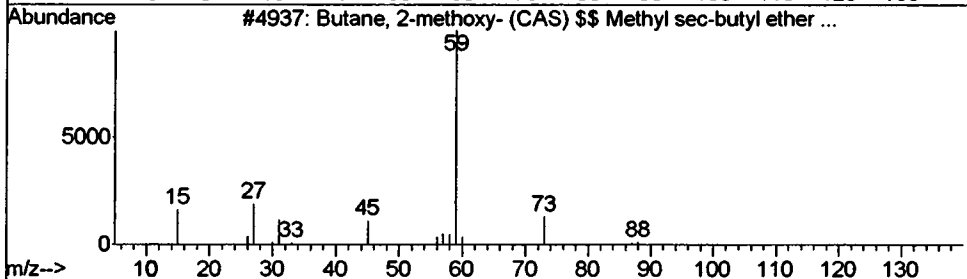
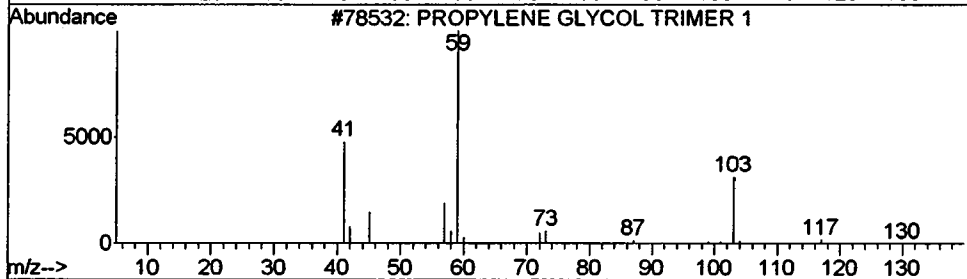
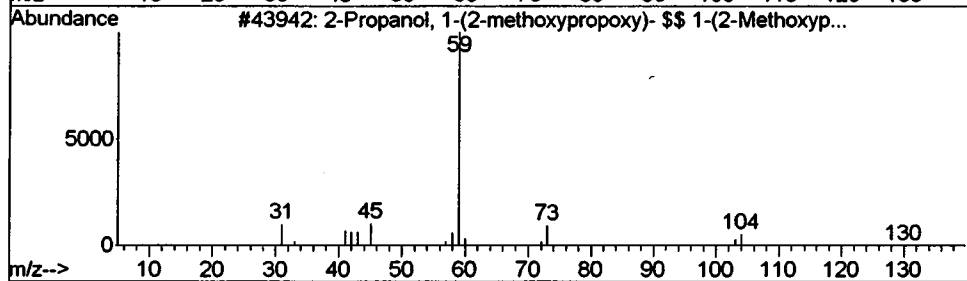
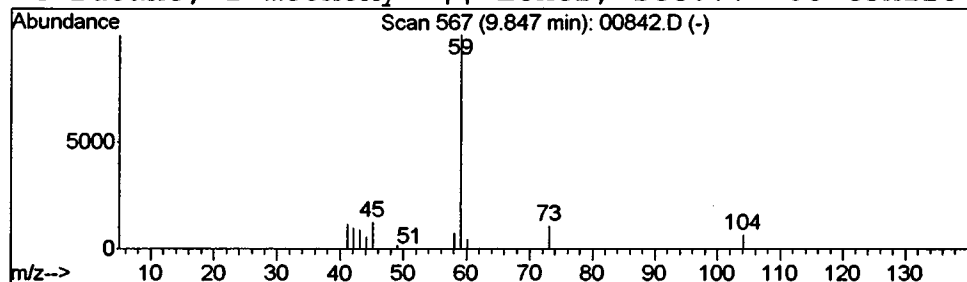
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Quant Method : C:\MSDCHEM\1...\HP020702.M (RTE Integrator)
 Title : 508 scan method
 Library : C:\DATABASE\WILEY7N.L

 Peak Number 8 2-Propanol, 1-(2-methoxypro... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.85	0.13 ug/L	52151	1,4-Dichlorobenzene-d4	9.71

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Propanol, 1-(2-methoxypropoxy)...	148	C7H16O3	013429-07-7	83
2			PROPYLENE GLYCOL TRIMER 1	174	C9H18O3	000000-00-0	64
3			Butane, 2-methoxy- (CAS) \$\$ Meth...	88	C5H12O	006795-87-5	40
4			Butane, 2-methoxy- \$\$ Ether, sec...	88	C5H12O	006795-87-5	40



Data File : C:\MSDCHEM\1\5973N\02FEB07\00842.D
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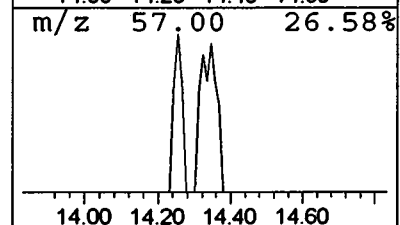
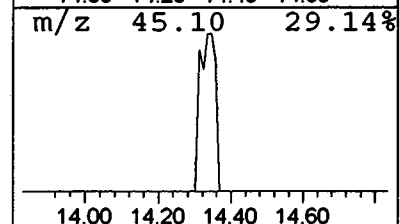
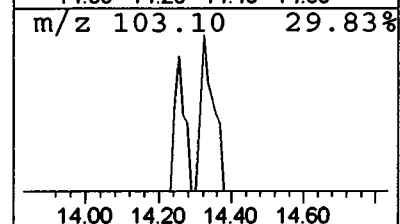
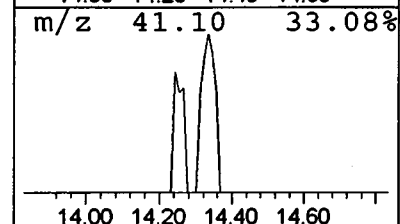
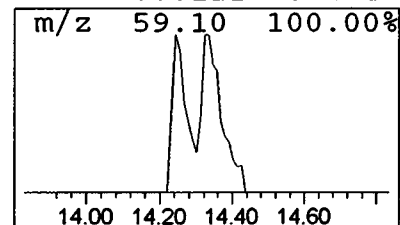
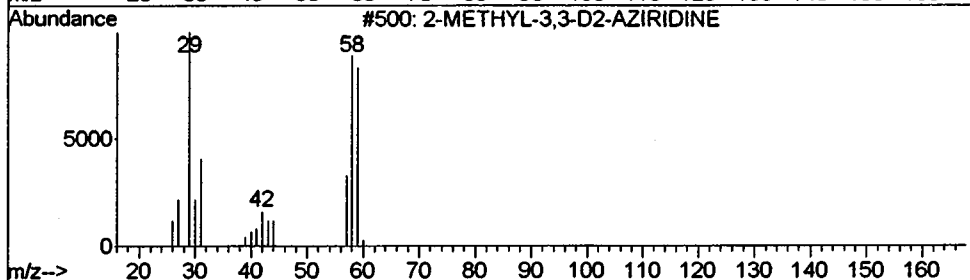
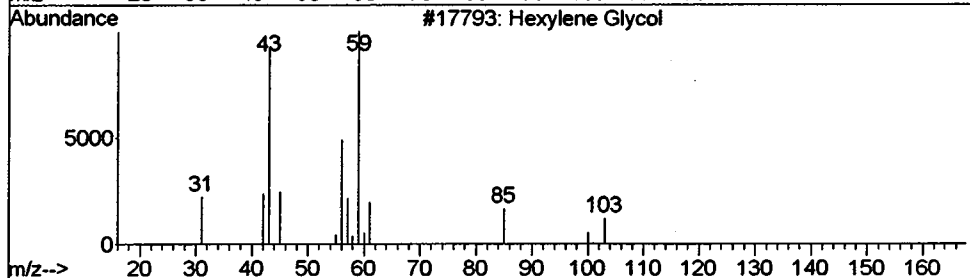
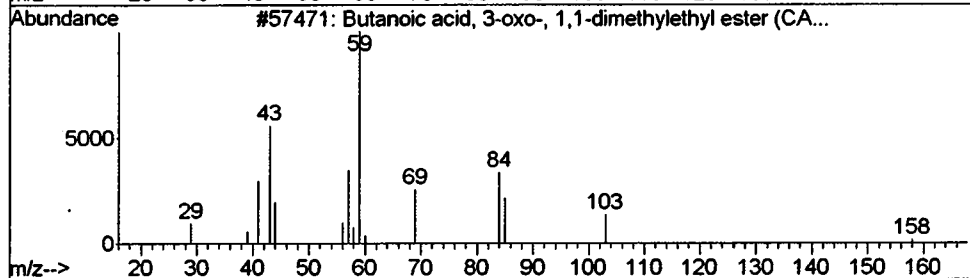
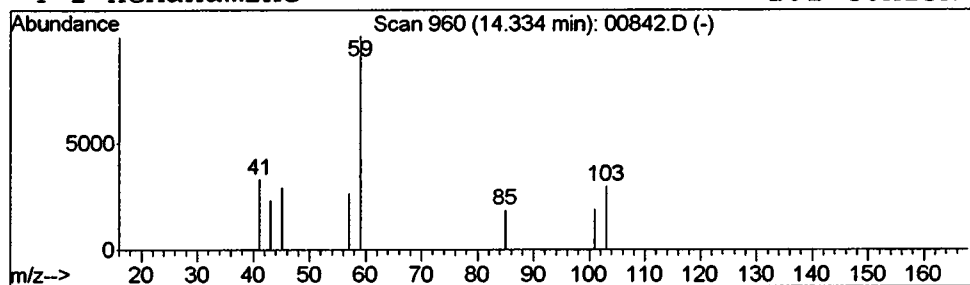
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 Multiplr: 1.00

Quant Method : C:\MSDCHEM\1...\HP020702.M (RTE Integrator)
 Title : 508 scan method
 Library : C:\DATABASE\WILEY7N.L

 Peak Number 9 Butanoic acid, 3-oxo-, 1,1-... Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.33	0.05 ug/L	31120	Naphthalene-d8	13.19

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Butanoic acid, 3-oxo-, 1,1-dimet...	158	C8H14O3	001694-31-1	9	
2	Hexylene Glycol	118	C6H14O2	000107-41-5	9	
3	2-METHYL-3,3-D2-AZIRIDINE	59	C3H5D2N	030691-55-5	4	
4	1-Hexanamine	101	C6H15N	000111-26-2	4	



Data File : C:\MSDCHEM\1\5973N\02FEB07\00842.D
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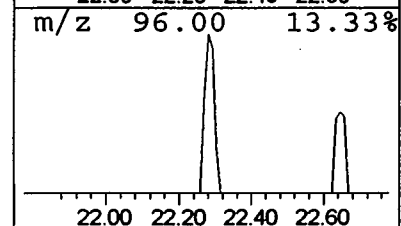
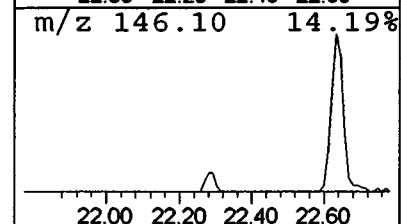
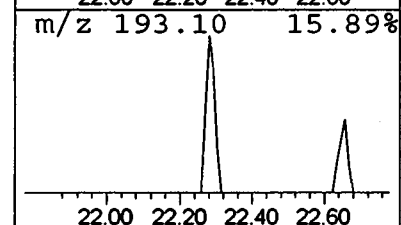
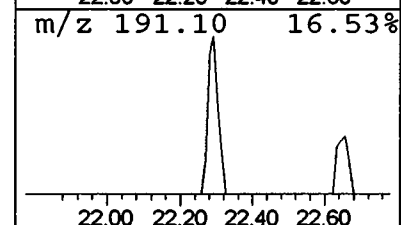
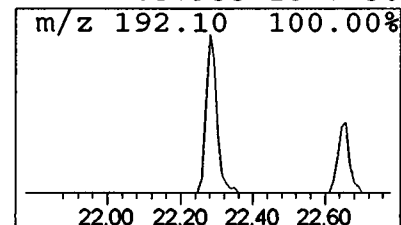
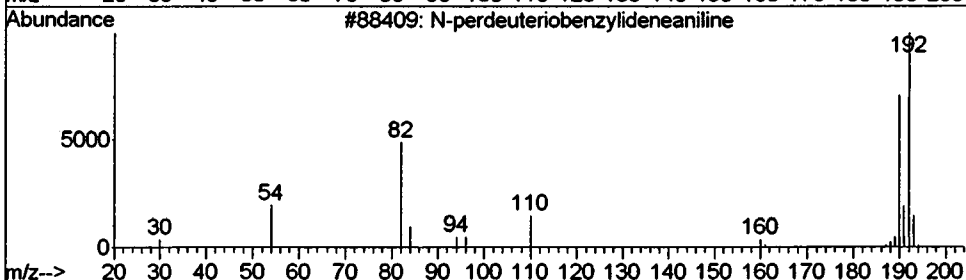
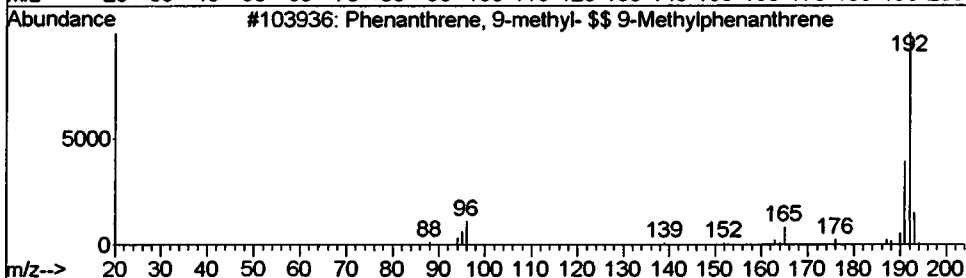
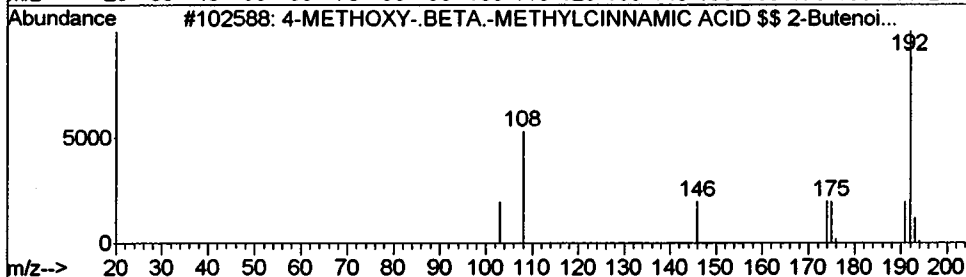
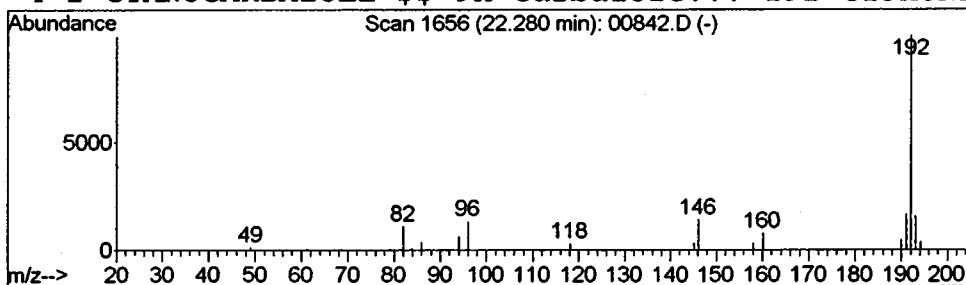
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Quant Method : C:\MSDCHEM\1...\HP020702.M (RTE Integrator)
 Title : 508 scan method
 Library : C:\DATABASE\WILEY7N.L

 Peak Number 10 4-METHOXY-.BETA.-METHYLCINN... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.28	0.13 ug/L	89734	Phenanthrene-d10	22.63

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			4-METHOXY-.BETA.-METHYLCINNAMIC ...	192	C11H12O3	019169-94-9	72
2			Phenanthrene, 9-methyl- \$\$ 9-Met...	192	C15H12	000883-20-5	59
3			N-perdeuteriobenzylideneaniline	192	C13D11N	000538-51-2	58
4			2-CYANOCARBAZOLE \$\$ 9H-Carbazole...	192	C13H8N2	057955-18-7	50



Data File : C:\MSDCHEM\1\5973N\02FEB07\00842.D
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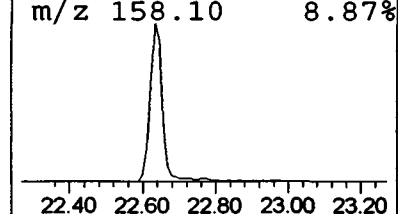
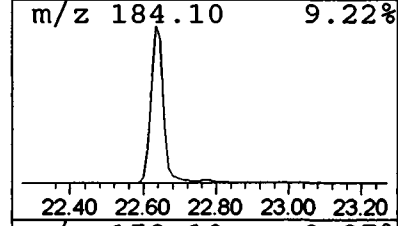
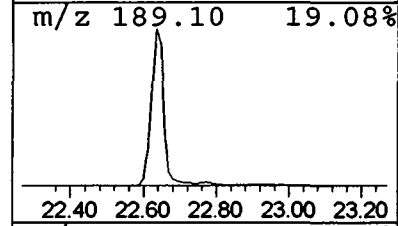
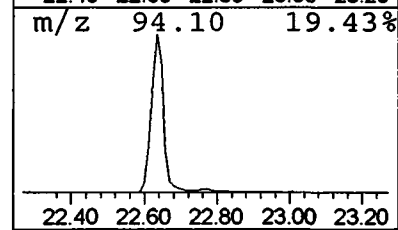
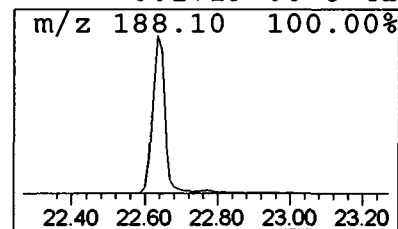
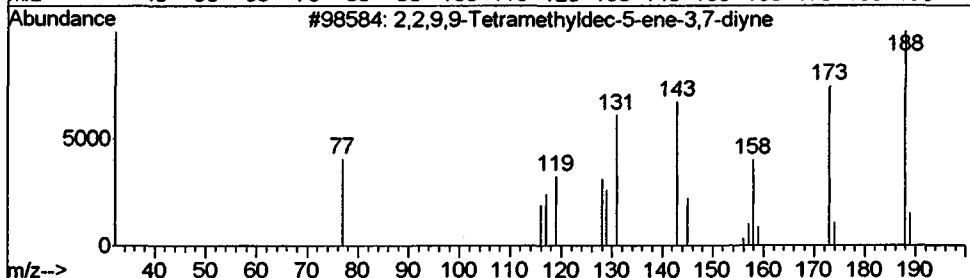
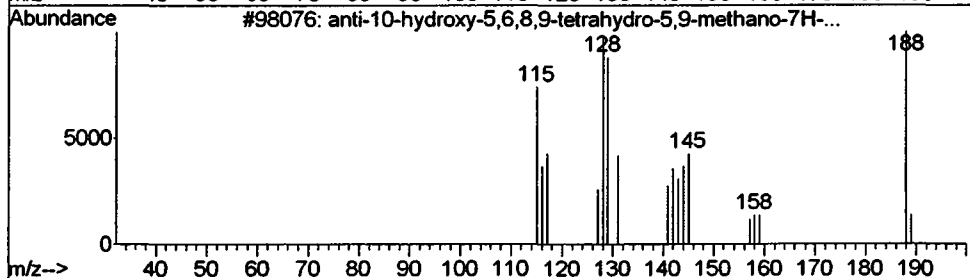
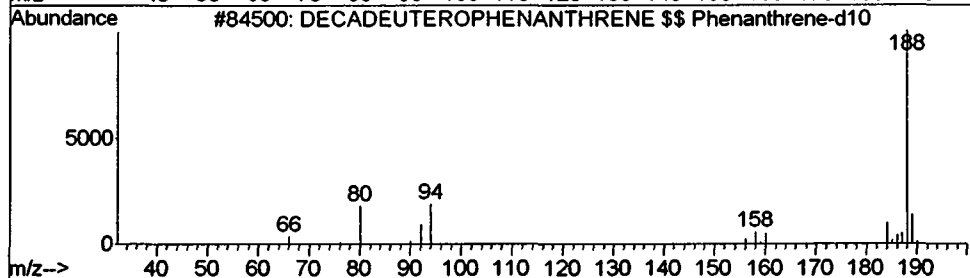
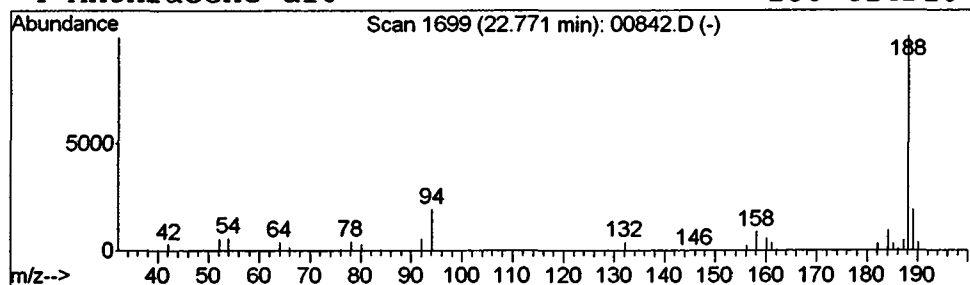
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Quant Method : C:\MSDCHEM\1...\HP020702.M (RTE Integrator)
 Title : 508 scan method
 Library : C:\DATABASE\WILEY7N.L

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.77	0.68 ug/L	451044	Phenanthrene-d10	22.63

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			DECADEUTEROPHENANTHRENE \$\$ Phena...	188	C14D10	001517-22-2	86
2			anti-10-hydroxy-5,6,8,9-tetrahyd...	188	C12H12O2	055139-53-2	58
3			2,2,9,9-Tetramethyldec-5-ene-3,7...	188	C14H20	102745-35-7	47
4			Anthracene-d10-	188	C14D10	001719-06-8	42



Data File : C:\MSDCHEM\1\5973N\02FEB07\00842.D
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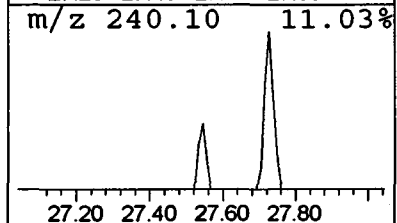
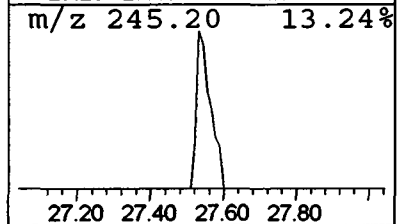
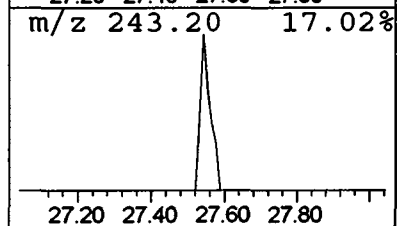
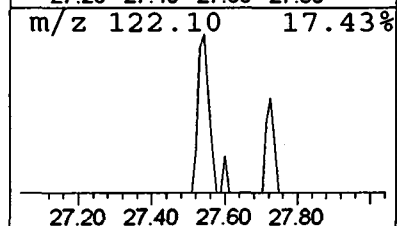
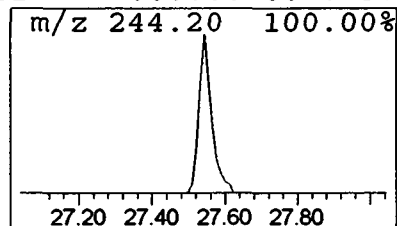
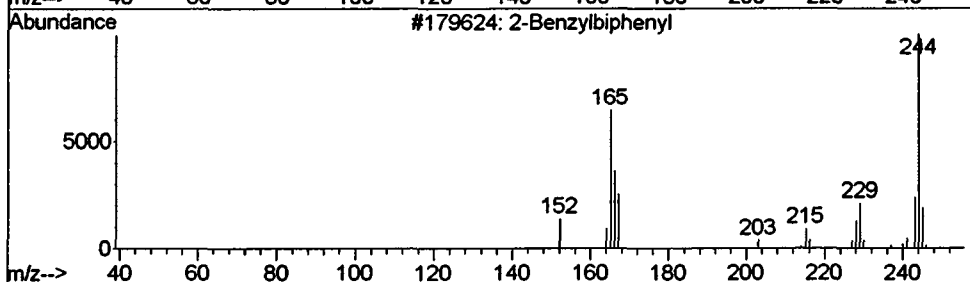
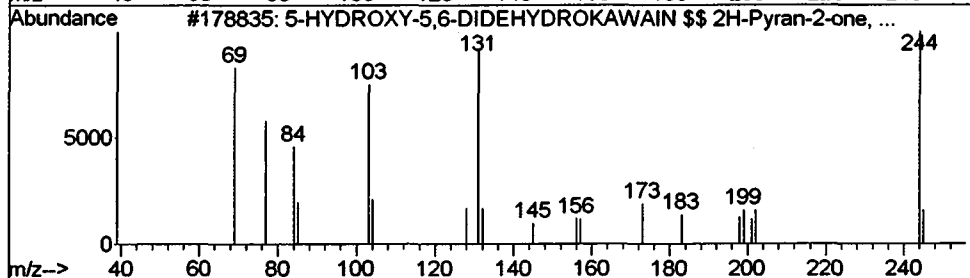
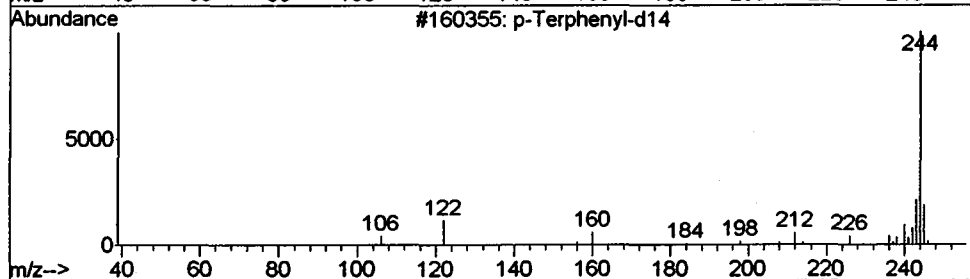
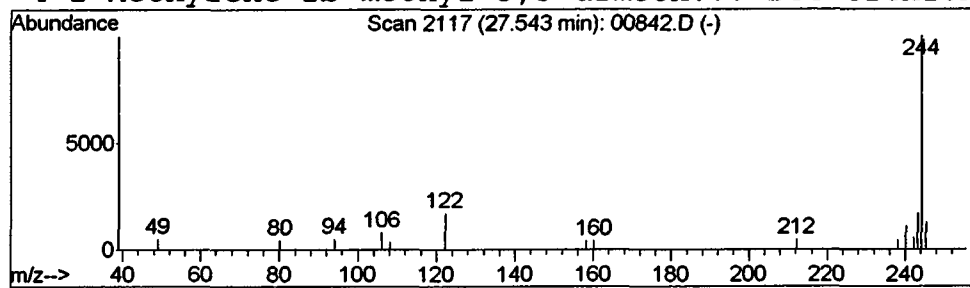
Vial: 6
 Operator:
 Inst : Instrumen
 Multiplr: 1.00

Quant Method : C:\MSDCHEM\1...\HP020702.M (RTE Integrator)
 Title : 508 scan method
 Library : C:\DATABASE\WILEY7N.L

 Peak Number 12 p-Terphenyl-d14 Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
27.54	0.11 ug/L	67550	Chrysene-d12	30.49

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			p-Terphenyl-d14	244	C18D14	001718-51-0	91
2			5-HYDROXY-5,6-DIDEHYDROKAWAIN \$\$...	244	C14H12O4	060037-36-7	58
3			2-Benzylbiphenyl	244	C19H16	000000-00-0	42
4			2-Methylene-2b-methyl-3,3-dimeth...	244	C14H16N2O2	000000-00-0	37



Data File : C:\MSDCHEM\1\5973N\02FEB07\00842.D
 Acq On : 7 Feb 2002 6:34 pm
 Sample : 508 scan
 Misc : February 7, 2002
 MS Integration Params: LSCINT.P

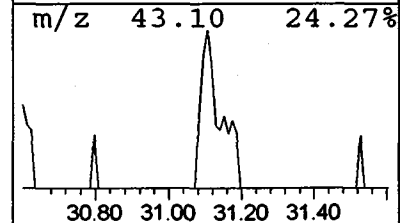
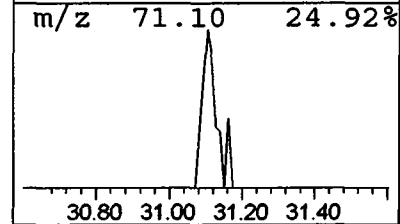
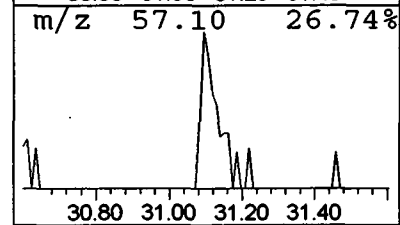
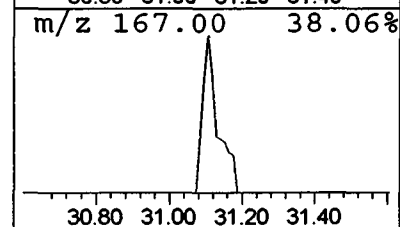
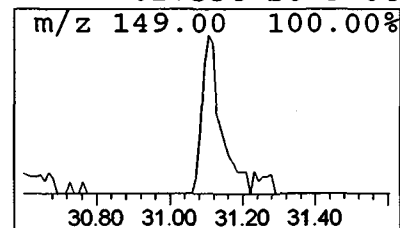
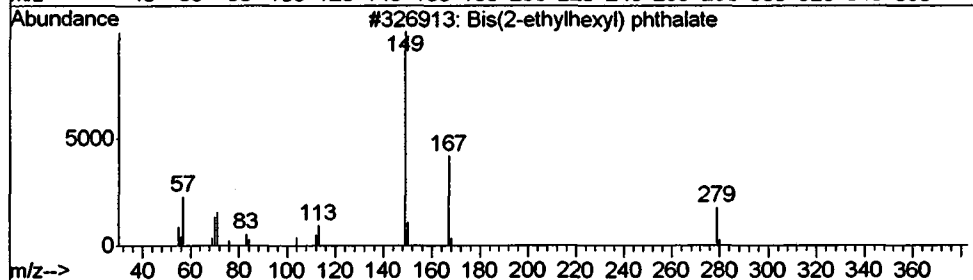
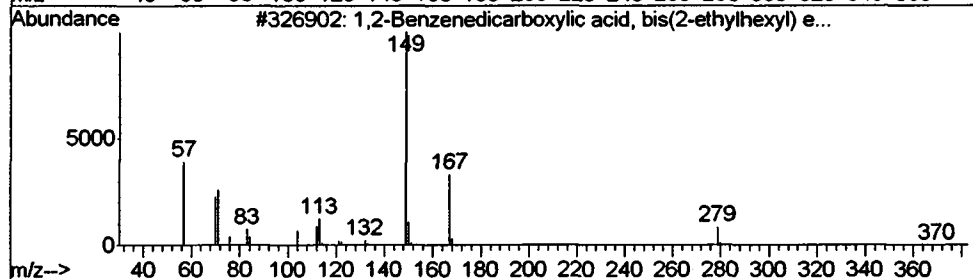
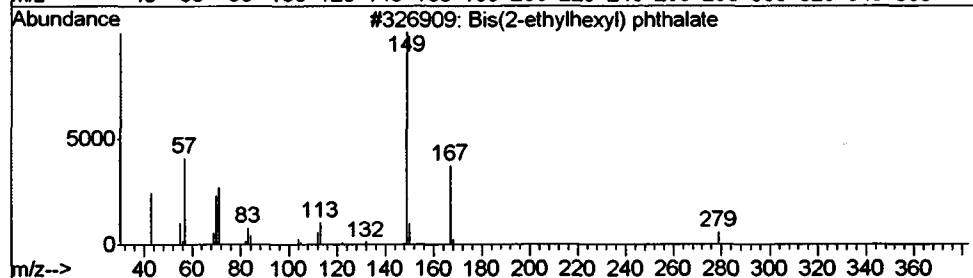
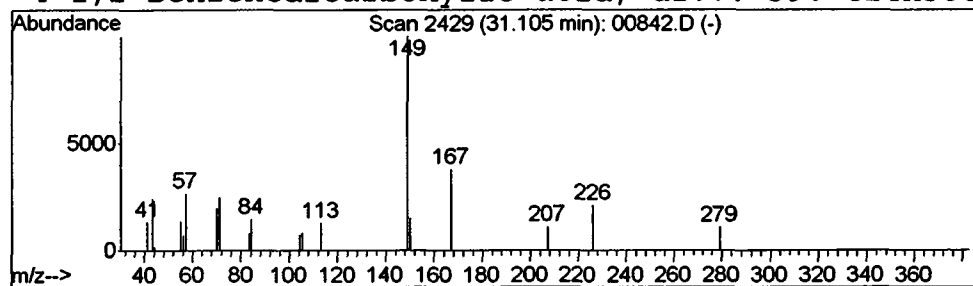
Vial: 6
 Operator:
 Inst : Instrumen
 Multiplr: 1.00

Quant Method : C:\MSDCHEM\1...\HP020702.M (RTE Integrator)
 Title : 508 scan method
 Library : C:\DATABASE\WILEY7N.L

 Peak Number 13 Bis(2-ethylhexyl) phthalate Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
31.11	0.19 ug/L	118534	Chrysene-d12	30.49

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Bis(2-ethylhexyl) phthalate	390	C24H38O4	000117-81-7	72
2			1,2-Benzenedicarboxylic acid, bi...	390	C24H38O4	000117-81-7	72
3			Bis(2-ethylhexyl) phthalate	390	C24H38O4	000117-81-7	64
4			1,2-Benzenedicarboxylic acid, di...	390	C24H38O4	027554-26-3	64



Operator ID: Date Acquired: 7 Feb 2002 6:34 pm
Data File: C:\MSDCHEM\1\5973N\02FEB07\00842.D
Name: 508 scan
Misc: February 7, 2002
Method: C:\MSDCHEM\1\METHODS\METHODS\HP020702.M (RTE Integrator)
Title: 508 scan method
Library Searched: C:\DATABASE\WILEY7N.L

TIC	Top Hit name	RT	EstConc	Units	Response	--Internal Standard---			
						#	RT	Resp	Conc
	Acetonitrile, dic...	3.49	0.2	ug/L	79931	1	9.71	8159580	20.0
	Butanoic acid, me...	3.61	0.1	ug/L	58894	1	9.71	8159580	20.0
	4-Penten-2-one	3.77	0.1	ug/L	39480	1	9.71	8159580	20.0
	Propane, 1,1-dime...	3.92	0.1	ug/L	32960	1	9.71	8159580	20.0
	3-(CYCLOHEX-3'-EN...	5.92	0.4	ug/L	172569	1	9.71	8159580	20.0
	2-Propanone, 1,1,...	6.01	0.4	ug/L	147393	1	9.71	8159580	20.0
	2-Propanol, 1-(2-...	9.56	0.1	ug/L	44022	1	9.71	8159580	20.0
	2-Propanol, 1-(2-...	9.85	0.1	ug/L	52151	1	9.71	8159580	20.0
	Butanoic acid, 3-...	14.33	0.1	ug/L	31120	2	13.19	11412400	20.0
	4-METHOXY-.BETA.-...	22.28	0.1	ug/L	89734	4	22.63	13319300	20.0
	DECADEUTEROPHENAN...	22.77	0.7	ug/L	451044	4	22.63	13319300	20.0
	p-Terphenyl-d14	27.54	0.1	ug/L	67550	5	30.49	12646300	20.0
	Bis(2-ethylhexyl)...	31.11	0.2	ug/L	118534	5	30.49	12646300	20.0