

User: Vinci, David L
Westchester Dept. of Labs & Research
Date: 02/04/2002 Time: 13:46:24

Sample: AE00843
Analysis: \$GCMS GCMS Scan For Unknowns
Units: ug
Started: 2/4/02 13:45 Ended: 2/4/02 13:45 Analyst: DLV
Page: 1

	Component Name	Viol	Result	Qualifier	MDL
1	Other unknown compounds		.		
2	Surr_2,4-DDD		82		5.0

Comments for Sample AE00843 - Analysis \$GCMS

Westchester Dept. of Labs & Research

User: Vinci, David L

Date: 02-04-2002 Time: 13:46:18

The GCMS scan, from 35 to 550 amu, does not reveal the presence of unusual compounds. The recoveries for 7 of the 43 compounds in the LCS Standard were above their acceptance ranges.

Quantitation Report (Not Reviewed)

Data File : C:\MSDCHEM\1\5973N\02JAN31\00843.D Vial: 5
 Acq On : 31 Jan 2002 6:01 pm Operator:
 Sample : 508 scan Inst : Instrumen
 Misc : January 31, 2002 Multiplr: 1.00
 MS Integration Params: SPLIT.P
 Quant Time: Feb 01 08:46:12 2002 Quant Results File: SCAN508.RES

Quant Method : C:\MSDCHEM\1...\SCAN508.M (RTE Integrator)
 Title : 508 scan method
 Last Update : Fri Jan 11 13:20:07 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	1929288	20.00	ug/L	0.00
10) Naphthalene-d8	13.19	136	8235587	20.00	ug/L	-0.02
17) Acenaphthene-d10	18.34	164	4323885	20.00	ug/L	-0.01
29) Phenanthrene-d10	22.65	188	6852705	20.00	ug/L	-0.02
38) Chrysene-d12	30.50	240	5765242	20.00	ug/L	-0.01
44) Perylene-d12	34.42	264	4156211	20.00	ug/L	0.00

System Monitoring Compounds		R.T.	QIon	Response	Conc	Units	Dev (Min)
37) 2,4-DDD		27.73	235	442722	4.09	ug/L	0.00
Spiked Amount	5.000			Recovery	=	81.80%	

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
20) Dimethylphthalate	18.34	163	702590	2.46	ug/L	# 56
21) 2,6-Dinitrotoluene	18.34	165	547525	9.83	ug/L	# 17
42) Bis(2-ethylhexyl)phthalate	31.11	149	47572	0.71	ug/L	# 87

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

00843.D SCAN508.M Fri Feb 01 08:46:13 2002

Page 1

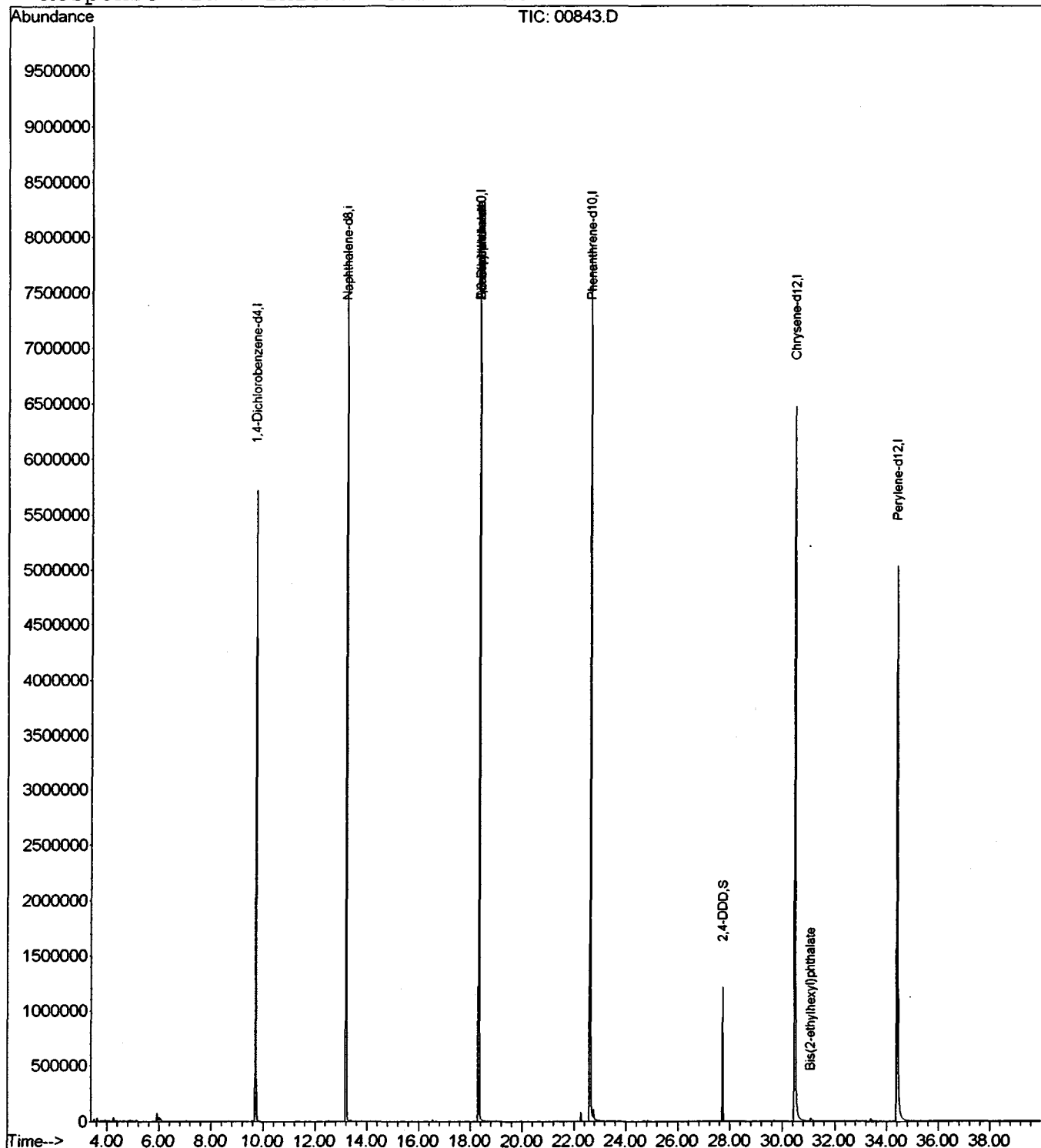
Quantitation Report (Not Reviewed)

Data File : C:\MSDCHEM\1\5973N\02JAN31\00843.D
 Acq On : 31 Jan 2002 6:01 pm
 Sample : 508 scan
 Misc : January 31, 2002
 MS Integration Params: SPLIT.P
 Quant Time: Feb 1 8:46 2002

Vial: 5
 Operator:
 Inst : Instrumen
 Multiplr: 1.00

Quant Results File: SCAN508.RE

Method : C:\MSDCHEM\1\METHODS\METHODS\SCAN508.M (RTE Integrator)
 Title : 508 scan method
 Last Update : Fri Jan 11 13:20:07 2002
 Response via : Initial Calibration



LSC Area Percent Report

Data File : C:\MSDCHEM\1\5973N\02JAN31\00843.D

Acq On : 31 Jan 2002 6:01 pm

Sample : 508 scan

Misc : January 31, 2002

MS Integration Params: LSCINT.P

Vial: 5

Operator:

Inst : Instrumen

Multiplr: 1.00

Method : C:\MSDCHEM\1\METHODS\METHODS\SCAN508.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

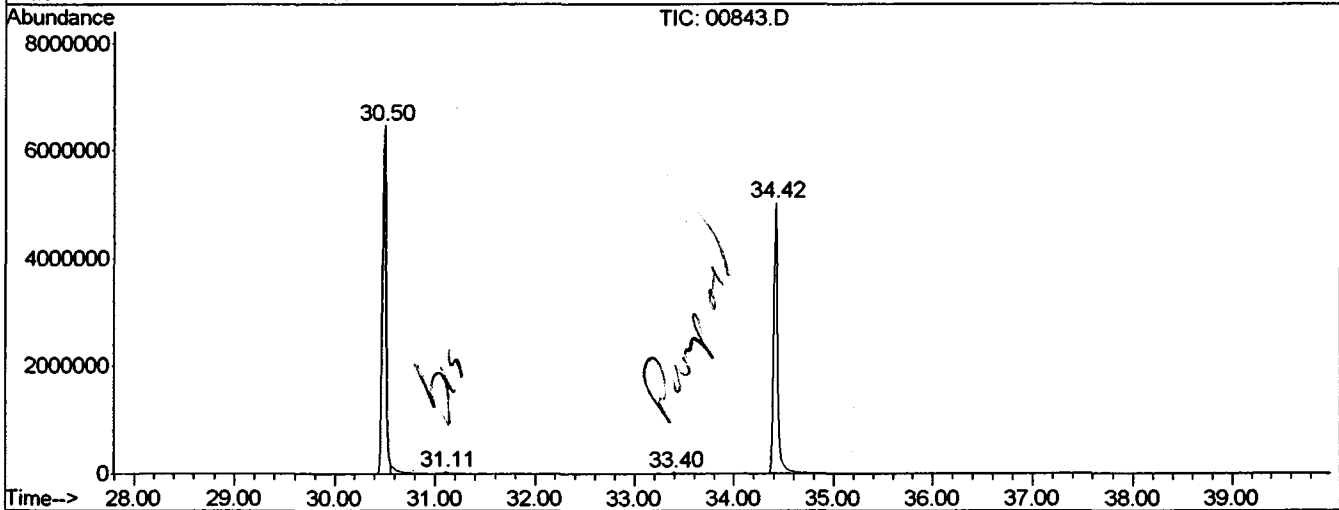
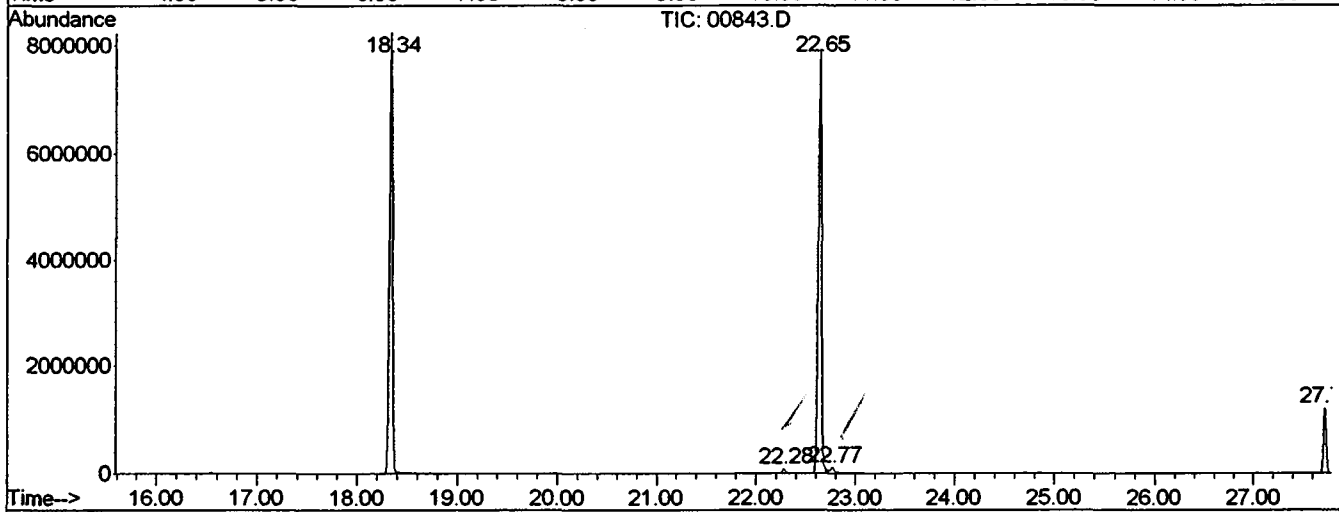
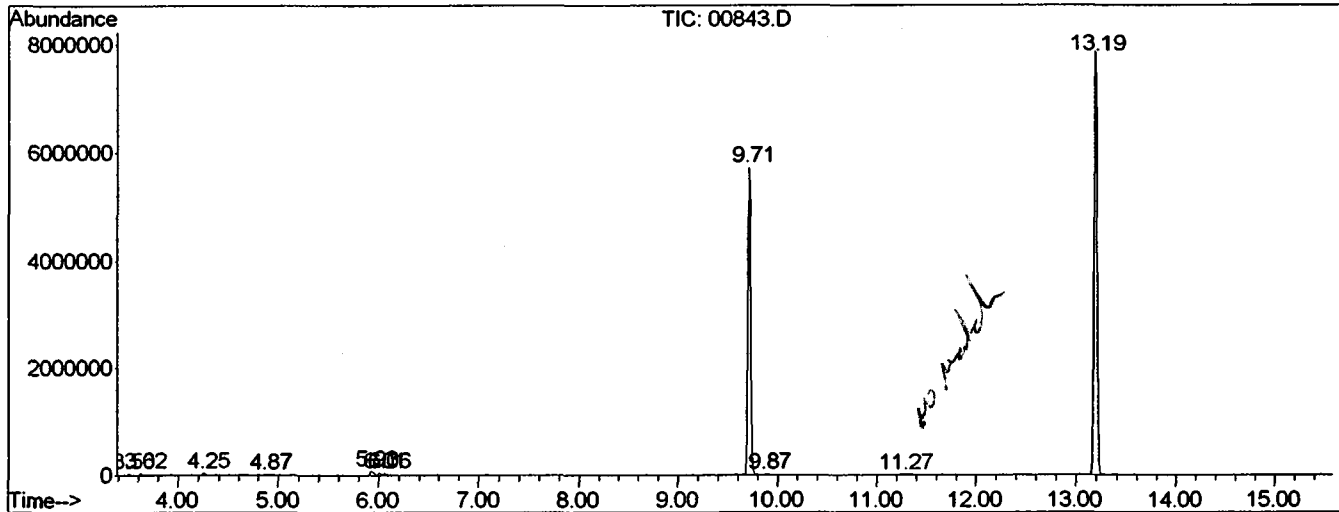
Signal : TIC

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.499	9	11	19	rVB	26893	45160	0.25%	0.046%
2	3.625	19	22	27	rBV	35032	60757	0.33%	0.062%
3	4.253	73	77	85	rVB3	38320	84914	0.47%	0.087%
4	4.869	129	131	141	rVV6	9323	42047	0.23%	0.043%
5	5.931	219	224	229	rBV	78008	189474	1.04%	0.195%
6	6.011	229	231	233	rVV	42609	71335	0.39%	0.073%
7	6.057	233	235	247	rVB	35802	91643	0.50%	0.094%
8	9.710	551	555	561	rVV	5720641	11147465	61.10%	11.450%
9	9.870	565	569	579	rVV	13411	47781	0.26%	0.049%
10	11.274	681	692	695	rBV4	9077	45385	0.25%	0.047%
11	13.192	855	860	865	rBV	7873397	16040535	87.92%	16.476%
12	18.341	1305	1311	1315	rBV	8250947	17429150	95.53%	17.903%
13	22.280	1651	1656	1661	rBV	86864	150114	0.82%	0.154%
14	22.646	1681	1688	1693	rBV2	7924568	18244770	100.00%	18.741%
15	22.771	1695	1699	1731	rVV	112506	443596	2.43%	0.456%
16	27.726	2129	2133	2137	rBV	1205299	2174277	11.92%	2.233%
17	30.500	2369	2376	2381	rBV	6466506	16668548	91.36%	17.121%
18	31.105	2423	2429	2443	rVV	34107	157887	0.87%	0.162%
19	33.400	2625	2630	2639	rBV	29264	94750	0.52%	0.097%
20	34.416	2713	2719	2763	rVB	5023103	14125094	77.42%	14.509%

Sum of corrected areas: 97354682

LSC Report - Integrated Chromatogram

File : C:\MSDCHEM\1\5973N\02JAN31\00843.D
 Operator :
 Acquired : 31 Jan 2002 6:01 pm using AcqMethod SCAN508
 Instrument : Instrumen
 Sample Name: 508 scan
 Misc Info : January 31, 2002
 Vial Number: 5
 Quant File :SCAN508.RES (RTE Integrator)



Library Search Compound Report

Data File : C:\MSDCHEM\1\5973N\02JAN31\00843.D
 Acq On : 31 Jan 2002 6:01 pm
 Sample : 508 scan
 Misc : January 31, 2002
 MS Integration Params: LSCINT.P

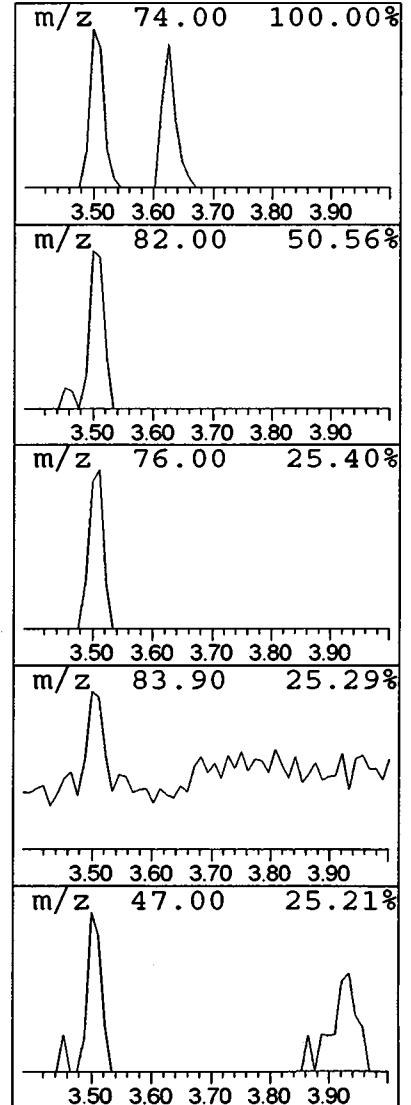
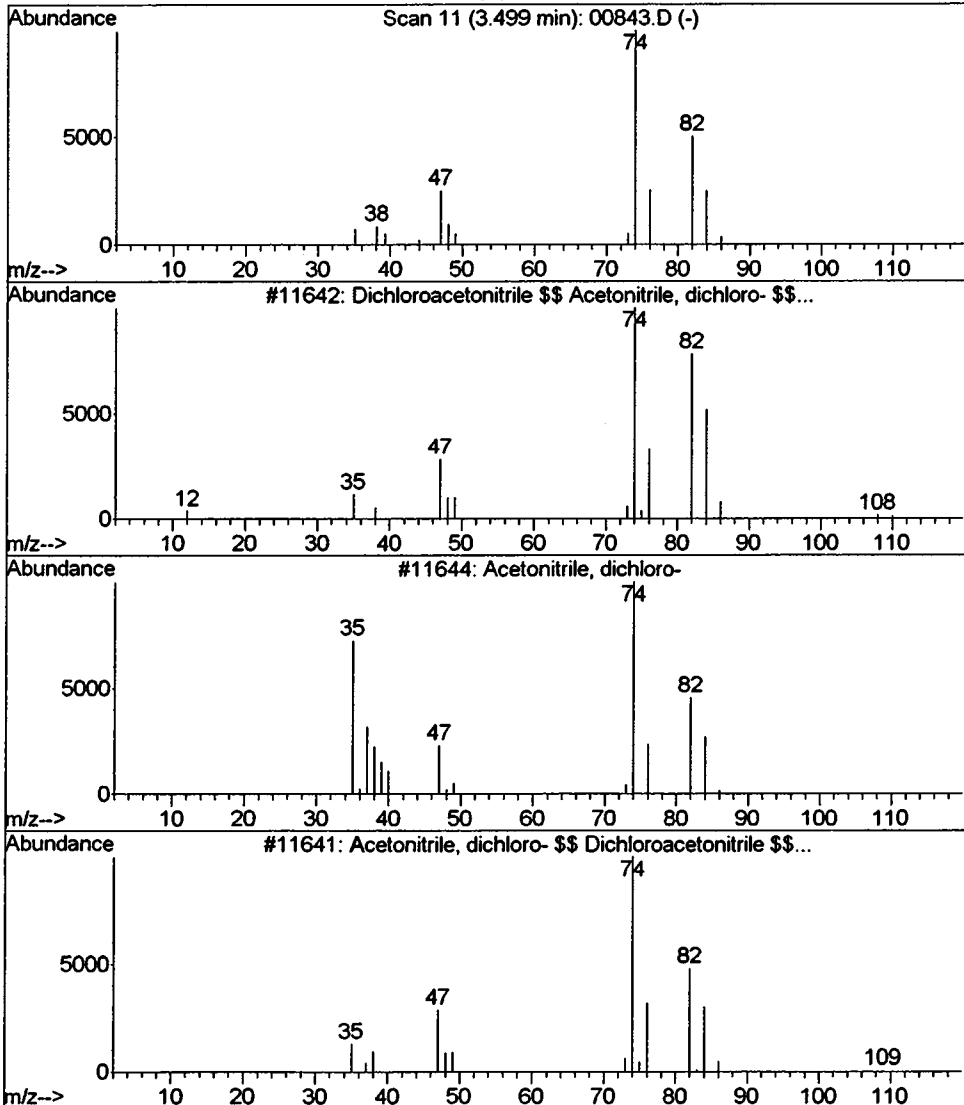
Vial: 5
 Operator:
 Inst : Instrumen
 Multiplr: 1.00

Quant Method : C:\MSDCHEM\1...\SCAN508.M (RTE Integrator)
 Title : 508 scan method
 Library : C:\DATABASE\WILEY7N.L

 Peak Number 1 Dichloroacetonitrile \$\$ Ace... Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.50	0.08 ug/L	45160	1,4-Dichlorobenzene-d4	9.71

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Dichloroacetonitrile \$\$ Acetonit...	109	C2HCl2N	003018-12-0	83
2			Acetonitrile, dichloro-	109	C2HCl2N	003018-12-0	64
3			Acetonitrile, dichloro- \$\$ Dichl...	109	C2HCl2N	003018-12-0	50
4			Dichloroacetonitrile \$\$ Acetonit...	109	C2HCl2N	003018-12-0	9



Library Search Compound Report

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 Acq On : 31 Jan 2002 6:01 pm
 Sample : 508 scan
 Misc : January 31, 2002
 MS Integration Params: LSCINT.P

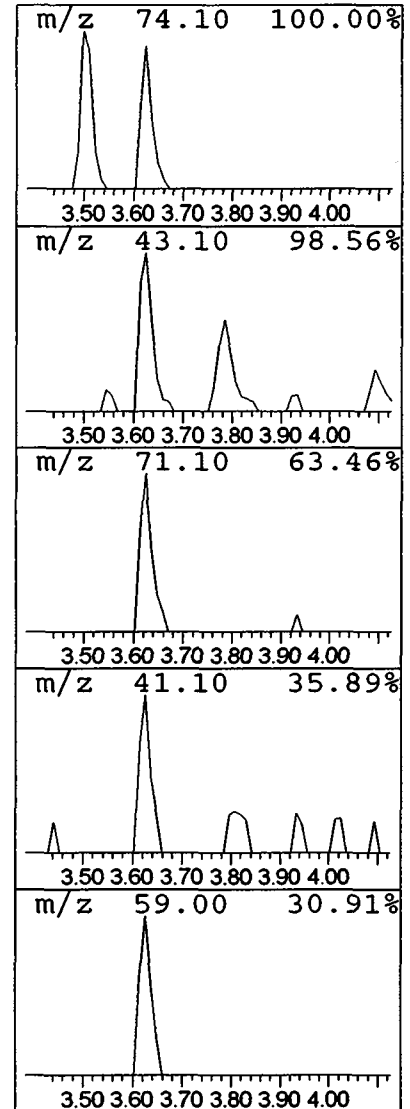
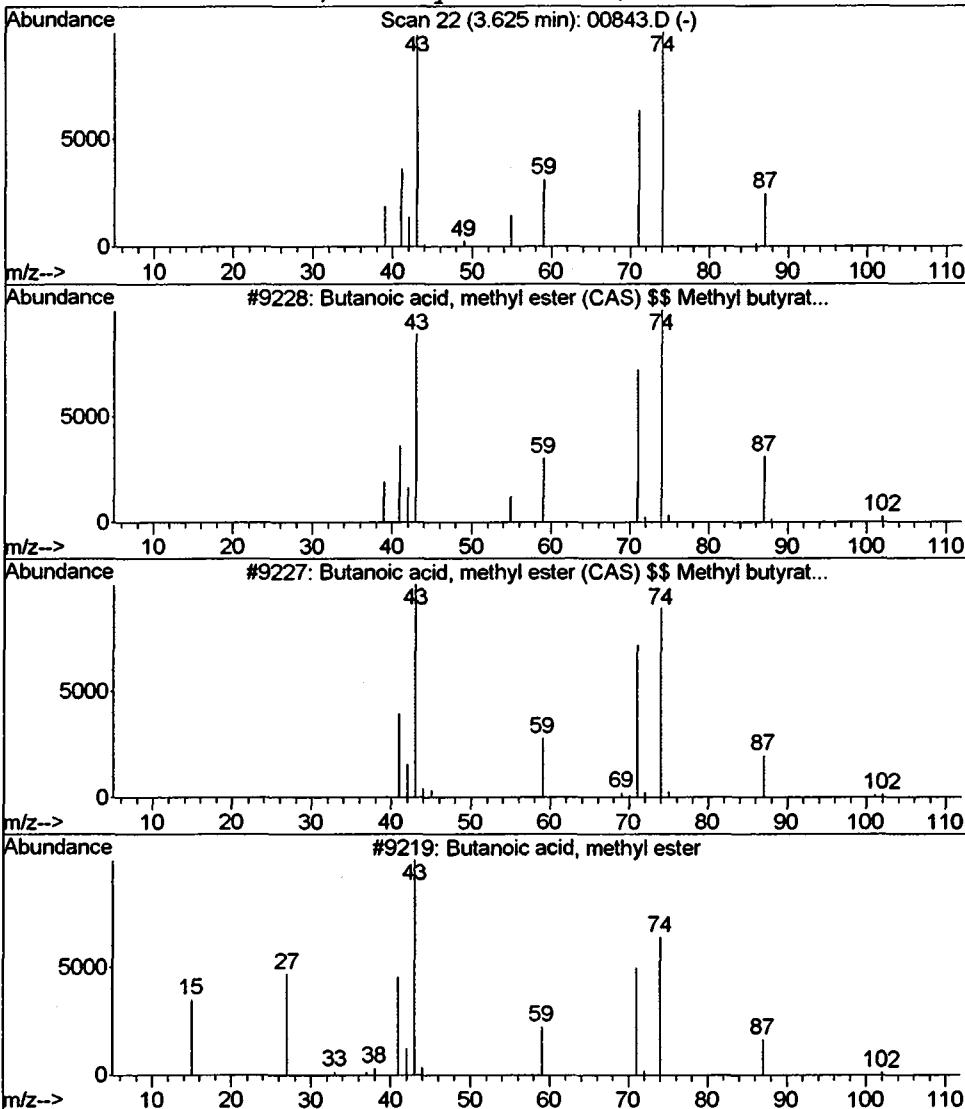
Vial: 5
 Operator:
 Inst : Instrumen
 Multiplr: 1.00

Quant Method : C:\MSDCHEM\1...\SCAN508.M (RTE Integrator)
 Title : 508 scan method
 Library : C:\DATABASE\WILEY7N.L

 Peak Number 2 Butanoic acid, methyl ester... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.62	0.11 ug/L	60757	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	102	C5H10O2	000623-42-7	83
4			Butanoic acid, methyl ester (CAS...	102	C5H10O2	000623-42-7	74



Library Search Compound Report

Data File : C:\MSDCHEM\1\5973N\02JAN31\00843.D
Acq On : 31 Jan 2002 6:01 pm
Sample : 508 scan
Misc : January 31, 2002
MS Integration Params: LSCINT.P

Vial: 5
Operator:
Inst : Instrumen
Multiplr: 1.00

Quant Method : C:\MSDCHEM\1...\SCAN508.M (RTE Integrator)

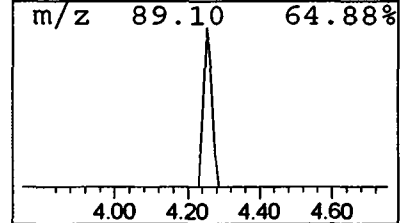
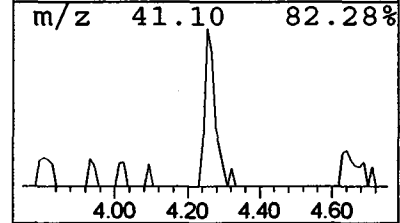
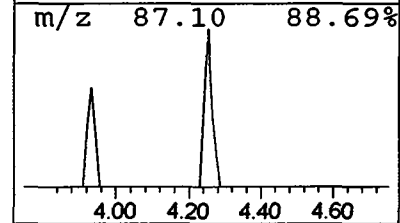
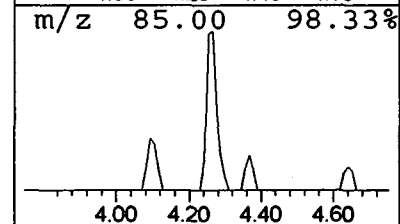
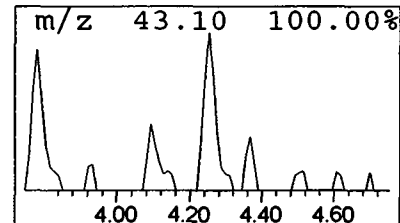
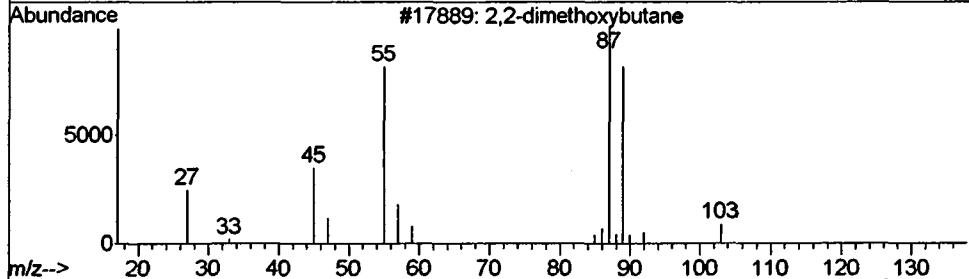
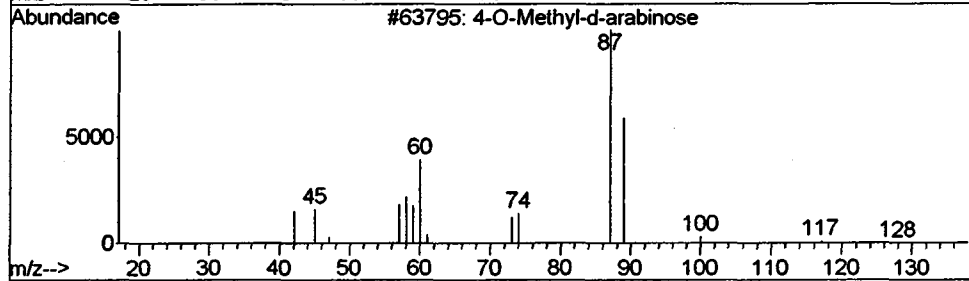
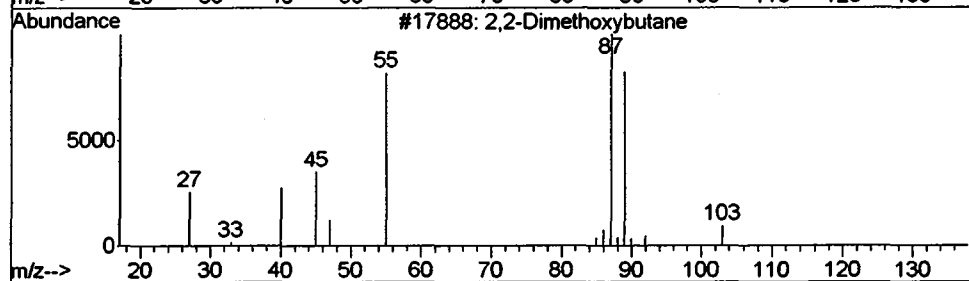
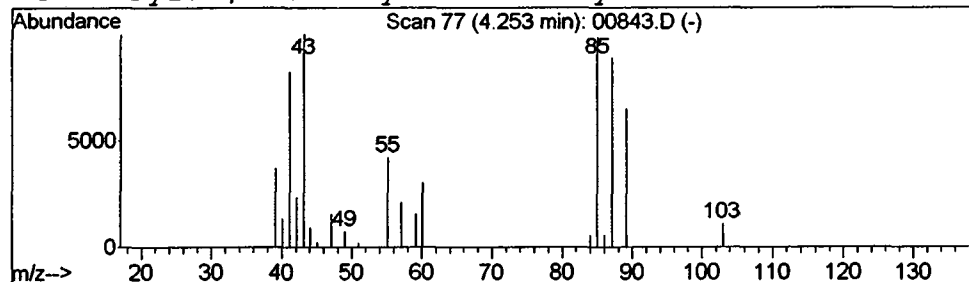
Title : 508 scan method

Library : C:\DATABASE\WILEY7N.L

Peak Number 3 2,2-Dimethoxybutane Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.25	0.15 ug/L	84914	1,4-Dichlorobenzene-d4	9.71

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2,2-Dimethoxybutane	118	C6H14O2	003453-99-4	38
2			4-O-Methyl-d-arabinose	164	C6H12O5	000000-00-0	37
3			2,2-dimethoxybutane	118	C6H14O2	003453-99-4	32
4			2H-Pyran, tetrahydro-2-methyl-	100	C6H12O	010141-72-7	27



Library Search Compound Report

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Sample : 508 scan
Misc : January 31, 2002
MS Integration Params: LSCINT.P

Vial: 5
Operator:
Inst : Instrumen
Multiplr: 1.00

Quant Method : C:\MSDCHEM\1...\SCAN508.M (RTE Integrator)

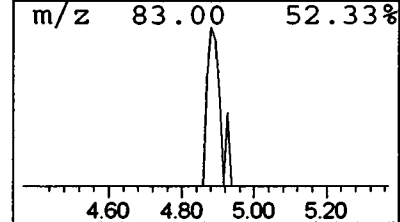
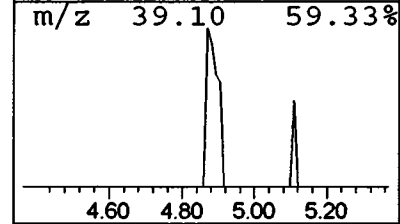
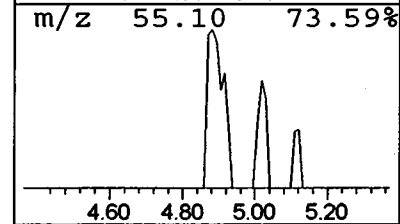
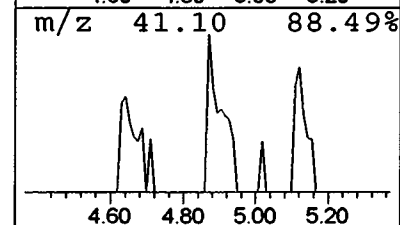
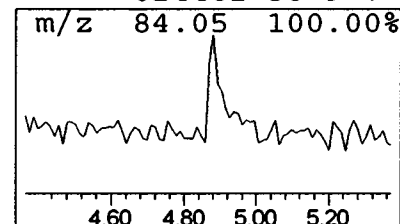
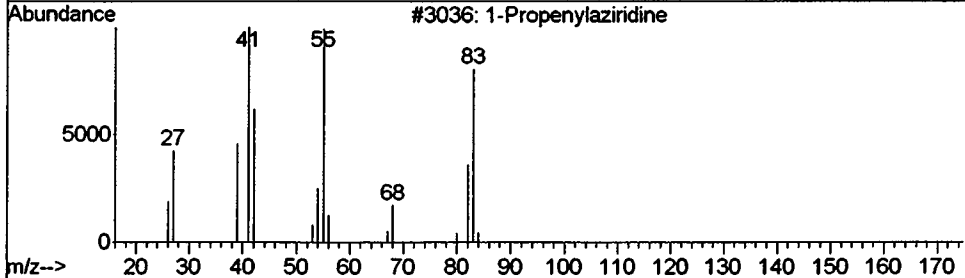
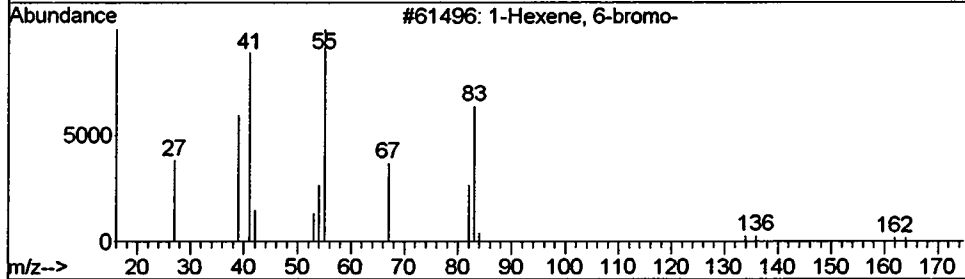
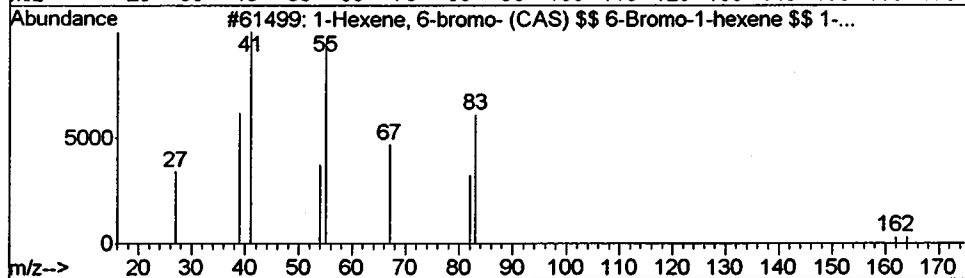
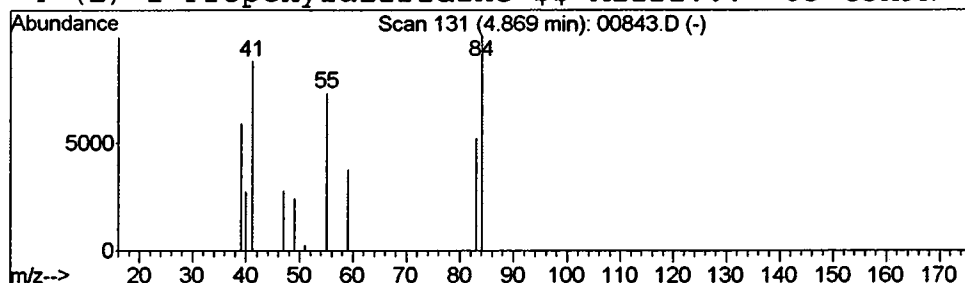
Title : 508 scan method

Library : C:\DATABASE\WILEY7N.L

Peak Number 4 1-Hexene, 6-bromo- (CAS) \$\$... Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.87	0.08 ug/L	42047	1,4-Dichlorobenzene-d4	9.71

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1-Hexene, 6-bromo- (CAS) \$\$ 6-Br...	162	C6H11Br	002695-47-8	9
2			1-Hexene, 6-bromo-	162	C6H11Br	002695-47-8	9
3			1-Propenylaziridine	83	C5H9N	000000-00-0	7
4			(Z)-1-Propenylaziridine \$\$ Aziri...	83	C5H9N	024461-38-9	7



Library Search Compound Report

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Vial: 5
 Operator:
 Inst : Instrumen
 Multiplr: 1.00

Quant Method : C:\MSDCHEM\1...\SCAN508.M (RTE Integrator)

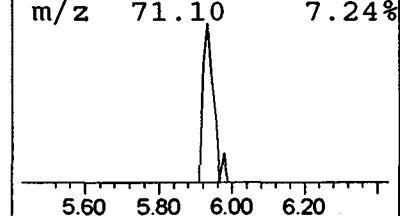
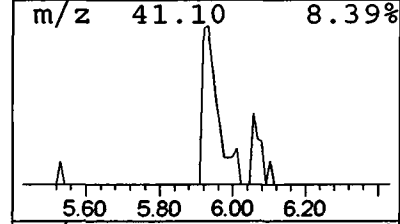
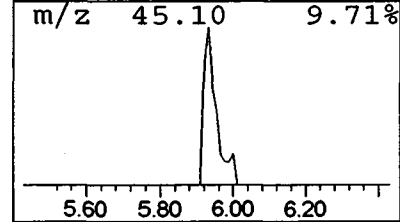
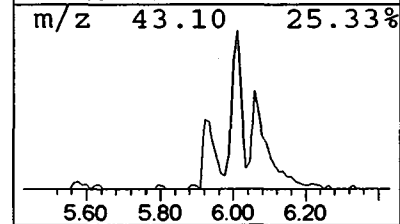
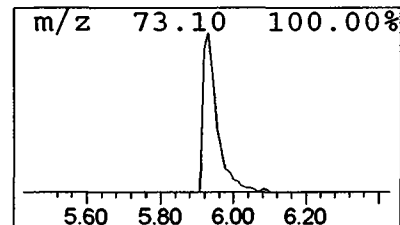
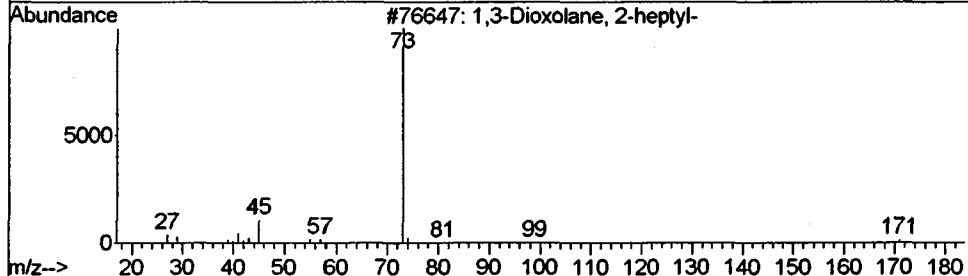
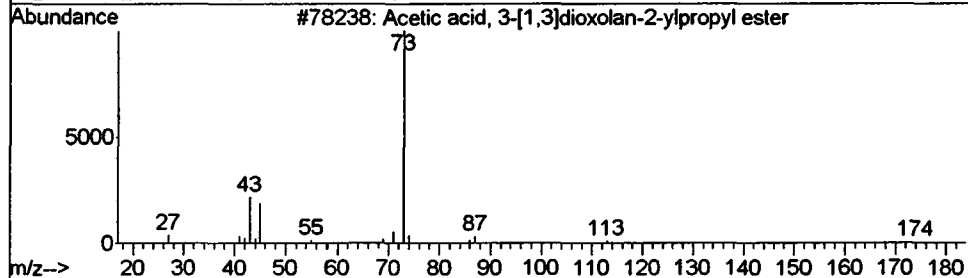
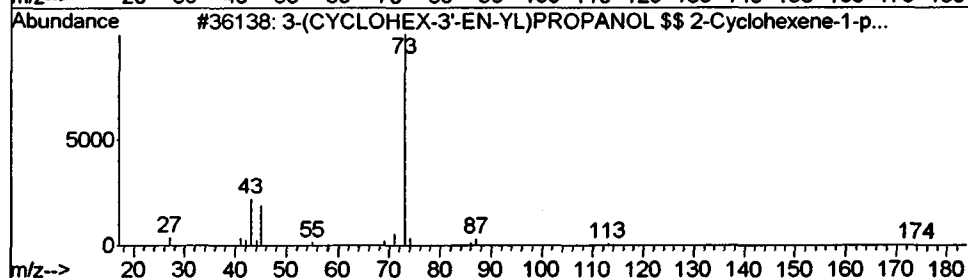
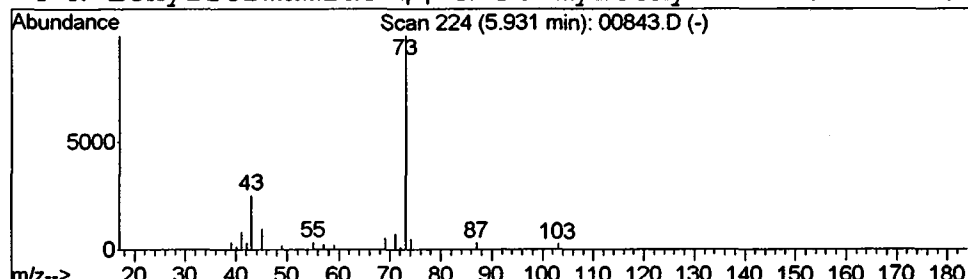
Title : 508 scan method

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.
5.93	0.34 ug/L	189474	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	28
2	Acetic acid, 3-[1,3]dioxolan-2-y...	174	C8H14O4	000000-00-0	28
3	1,3-Dioxolane, 2-heptyl-	172	C10H20O2	004359-57-3	25
4	N-Ethylformamide \$\$ N-Formylethy...	73	C3H7NO	000627-45-2	9



Library Search Compound Report

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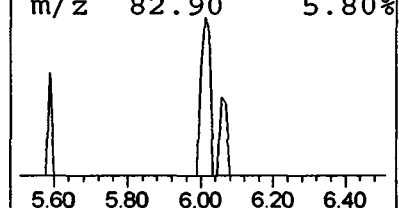
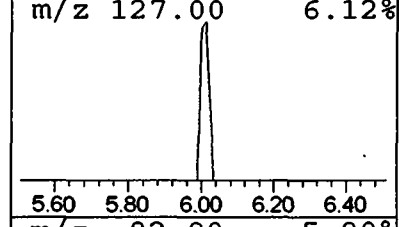
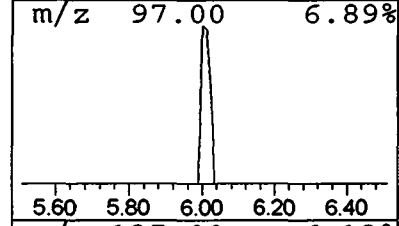
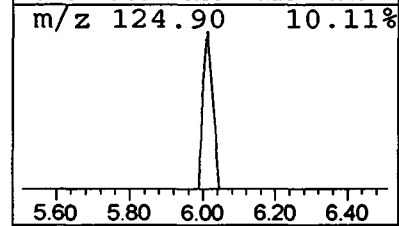
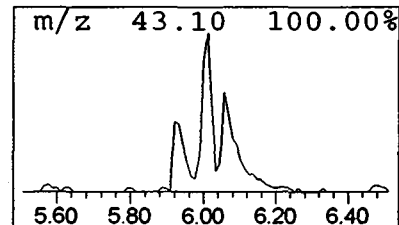
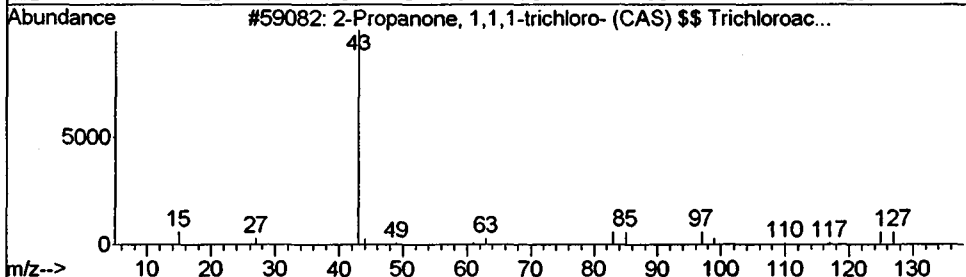
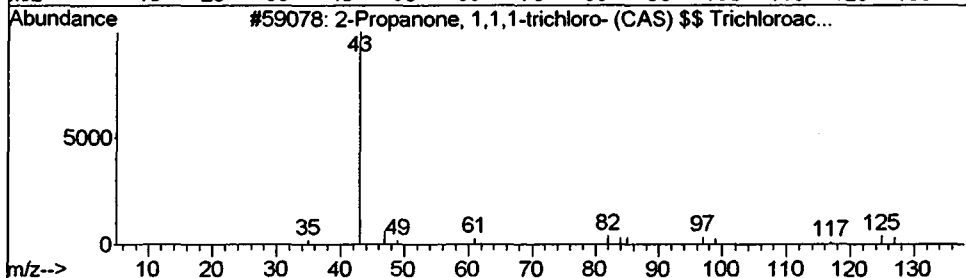
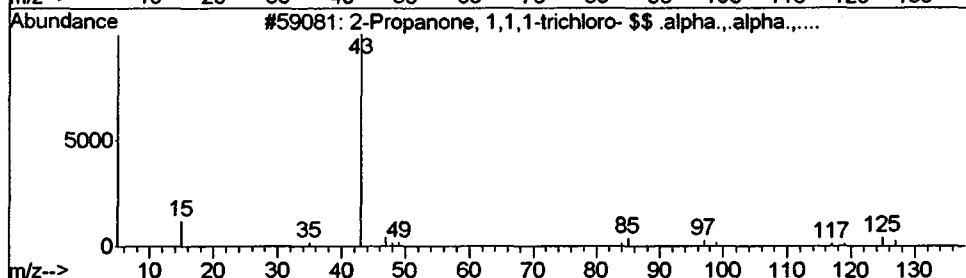
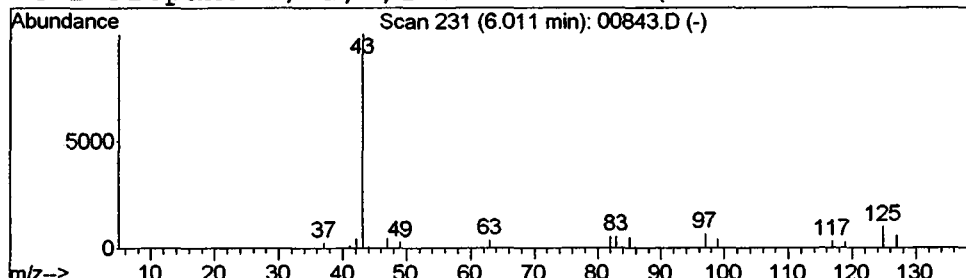
Vial: 5
 Operator:
 Inst : Instrumen
 Multiplr: 1.00

Quant Method : C:\MSDCHEM\1...\SCAN508.M (RTE Integrator)
 Title : 508 scan method
 Library : C:\DATABASE\WILEY7N.L

 Peak Number 6 2-Propanone, 1,1,1-trichlor... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.01	0.13 ug/L	71335	1,4-Dichlorobenzene-d4	9.71

Hit#	of	5	Tentative ID		MW	MolForm	CAS#	Qual
1	2-Propanone,	1,1,1-trichloro-	\$\$...	160	C3H3Cl3O	000918-00-3	78	
2	2-Propanone,	1,1,1-trichloro-	(C...	160	C3H3Cl3O	000918-00-3	50	
3	2-Propanone,	1,1,1-trichloro-	(C...	160	C3H3Cl3O	000918-00-3	45	
4	2-Propanone,	1,1,1-trichloro-	(C...	160	C3H3Cl3O	000918-00-3	45	



Library Search Compound Report

Data File : C:\MSDCHEM\1\5973N\02JAN31\00843.D
 Acq On : 31 Jan 2002 6:01 pm
 Sample : 508 scan
 Misc : January 31, 2002
 MS Integration Params: LSCINT.P

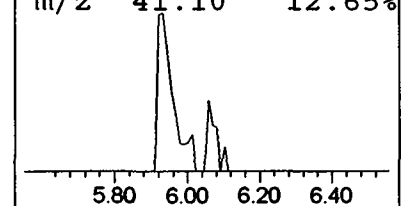
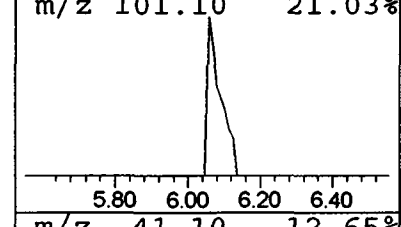
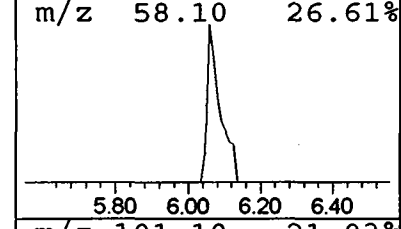
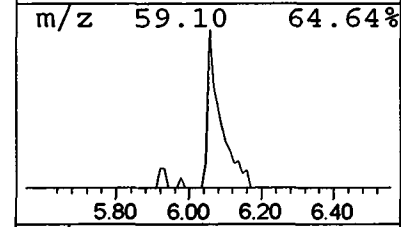
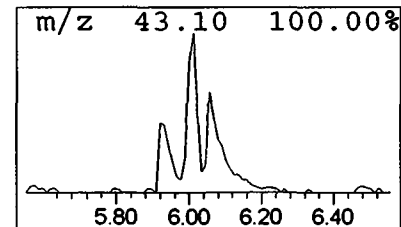
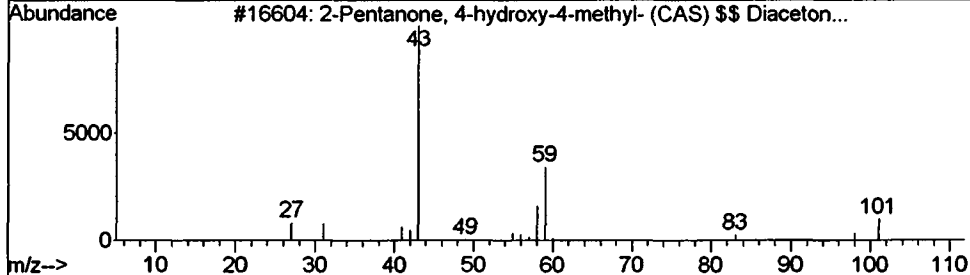
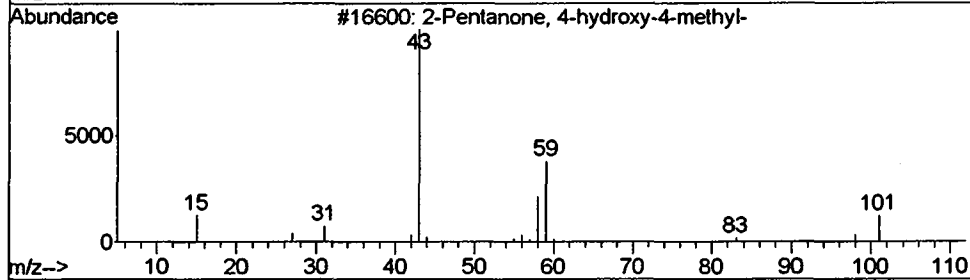
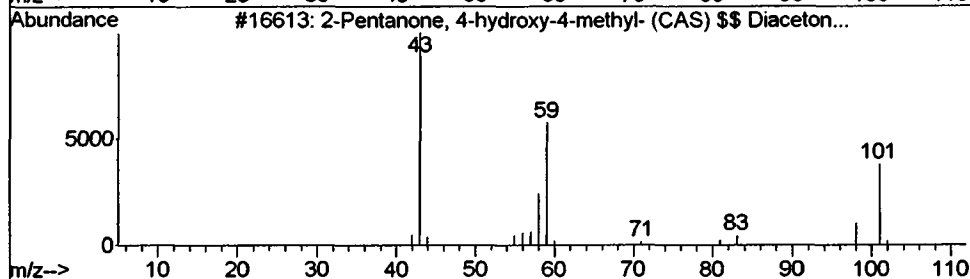
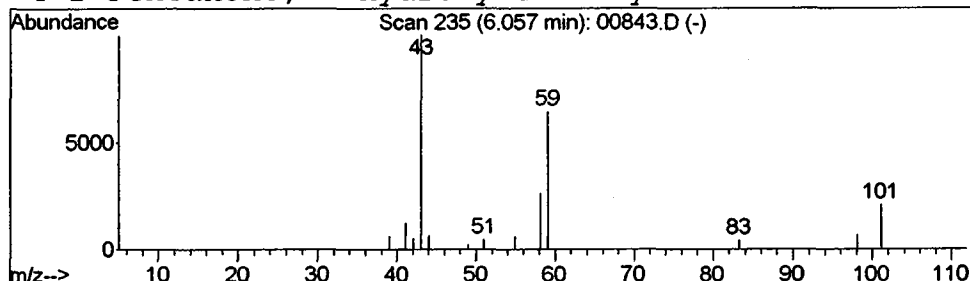
Vial: 5
 Operator:
 Inst : Instrumen
 Multiplr: 1.00

Quant Method : C:\MSDCHEM\1...\SCAN508.M (RTE Integrator)
 Title : 508 scan method
 Library : C:\DATABASE\WILEY7N.L

 Peak Number 7 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.06	0.16 ug/L	91643	1,4-Dichlorobenzene-d4	9.71

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentanone, 4-hydroxy-4-methyl-...	116	C6H12O2	000123-42-2	83
2			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	83
3			2-Pentanone, 4-hydroxy-4-methyl-...	116	C6H12O2	000123-42-2	78
4			2-Pentanone, 4-hydroxy-4-methyl-...	116	C6H12O2	000123-42-2	74



Library Search Compound Report

Data File : C:\MSDCHEM\1\5973N\02JAN31\00843.D
 Acq On : 31 Jan 2002 6:01 pm
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 Misc : January 31, 2002
 MS Integration Params: LSCINT.P

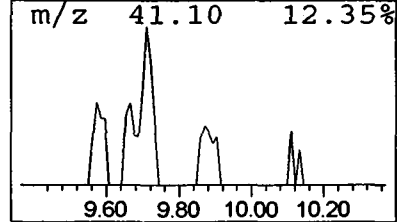
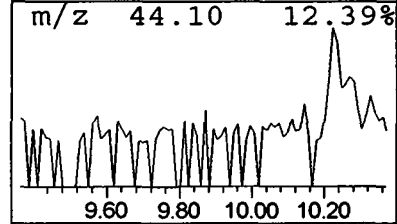
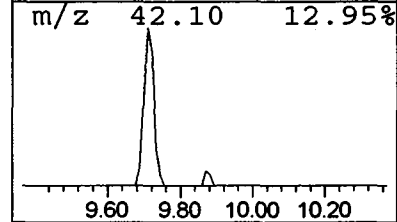
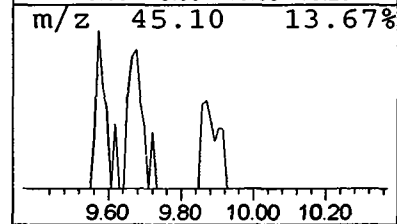
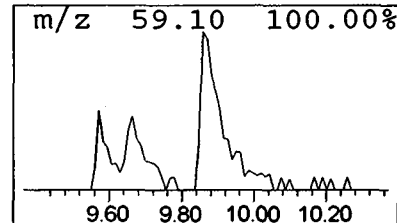
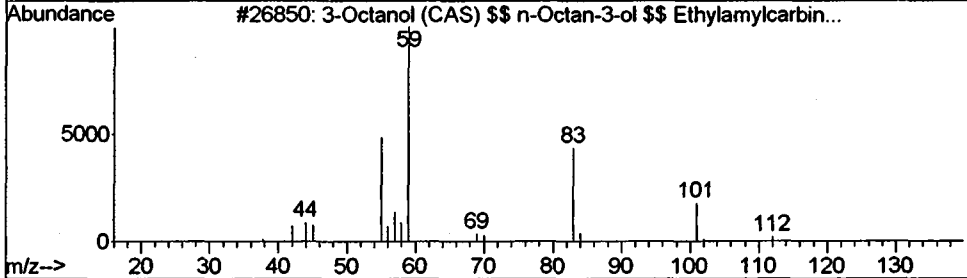
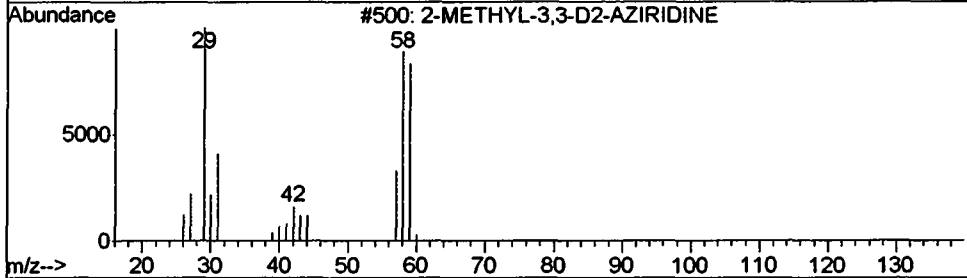
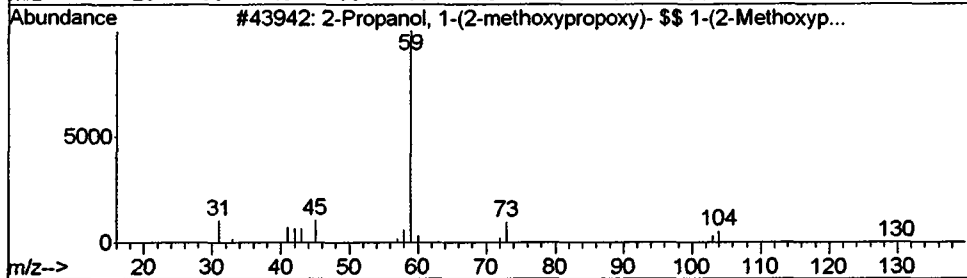
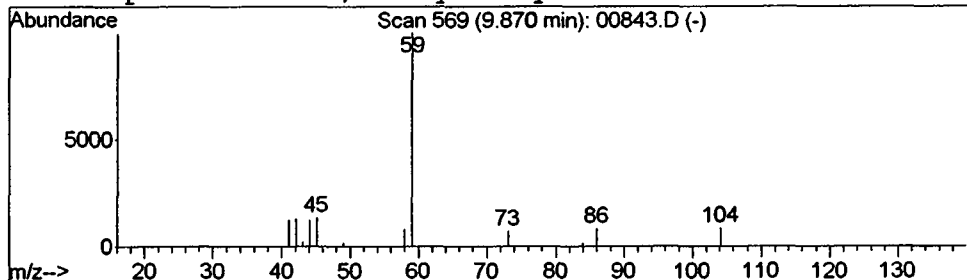
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 Inst : Instrumen
 Multiplr: 1.00

Quant Method : C:\MSDCHEM\1...\SCAN508.M (RTE Integrator)
 Title : 508 scan method
 Library : C:\DATABASE\WILEY7N.L

 Peak Number 8 2-Propanol, 1-(2-methoxypro... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.87	0.09 ug/L	47781	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	39
2		2-METHYL-3,3-D2-AZIRIDINE	59	C3H5D2N	030691-55-5	9
3		3-Octanol (CAS) \$\$ n-Octan-3-ol ...	130	C8H18O	000589-98-0	9
4		Propanoic acid, 2-hydroxy-2-meth...	118	C5H10O3	002110-78-3	9



Library Search Compound Report

Data File : C:\MSDCHEM\1\5973N\02JAN31\00843.D
 Acq On : 31 Jan 2002 6:01 pm
 Sample : 508 scan
 Misc : January 31, 2002
 MS Integration Params: LSCINT.P

Vial: 5
 Operator:
 Inst : Instrumen
 Multiplr: 1.00

Quant Method : C:\MSDCHEM\1...\SCAN508.M (RTE Integrator)

Title : 508 scan method

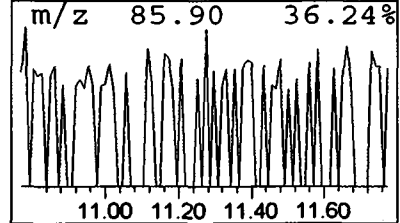
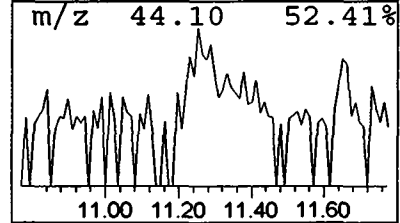
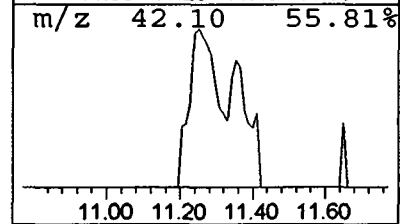
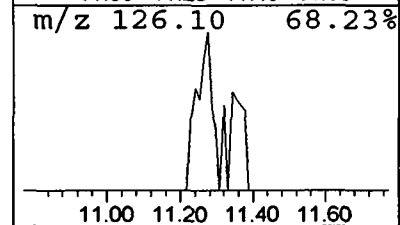
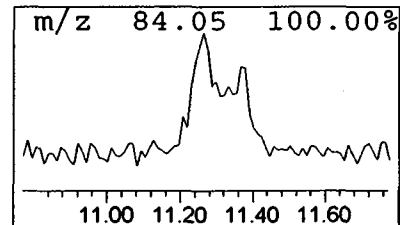
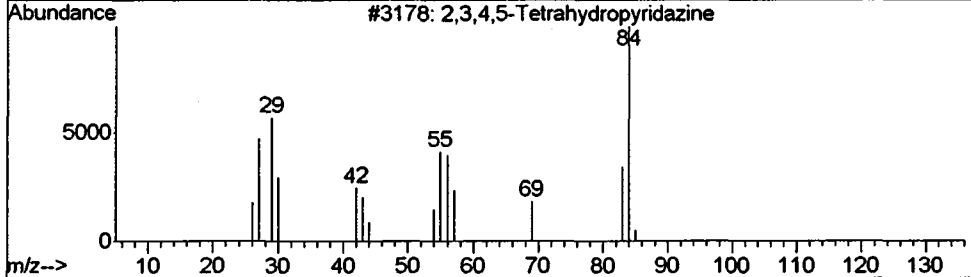
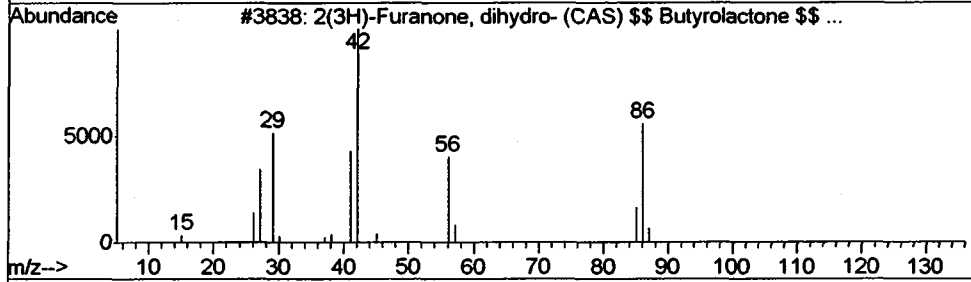
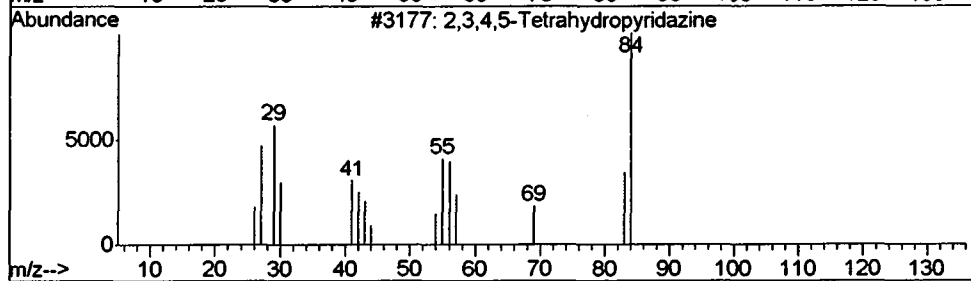
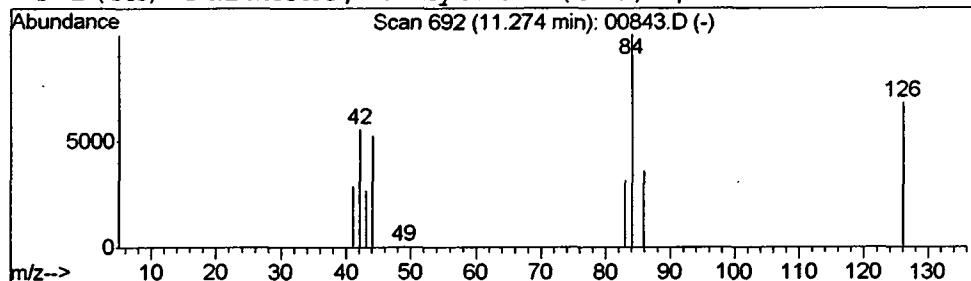
Library : C:\DATABASE\WILEY7N.L

 Peak Number 9 2,3,4,5-Tetrahydropyridazine Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.27	0.08 ug/L	45385	1,4-Dichlorobenzene-d4	9.71

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2,3,4,5-Tetrahydropyridazine	84	C4H8N2	000694-06-4	9
2			2(3H)-Furanone, dihydro- (CAS) \$...	86	C4H6O2	000096-48-0	9
3			2,3,4,5-Tetrahydropyridazine	84	C4H8N2	000694-06-4	7
4			2(3H)-Furanone, dihydro- (CAS) \$...	86	C4H6O2	000096-48-0	7

*NO
GOOD
MATCH*



Library Search Compound Report

Data File : C:\MSDCHEM\1\5973N\02JAN31\00843.D
Acq On : 31 Jan 2002 6:01 pm
Sample : 508 scan
Misc : January 31, 2002
MS Integration Params: LSCINT.P

Vial: 5
Operator:
Inst : Instrumen
Multiplr: 1.00

Quant Method : C:\MSDCHEM\1...\SCAN508.M (RTE Integrator)

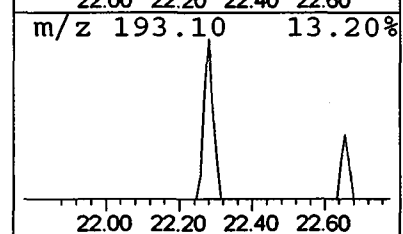
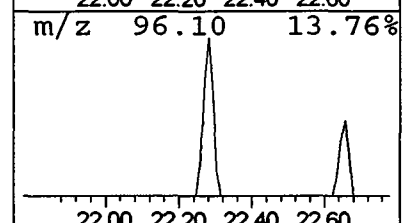
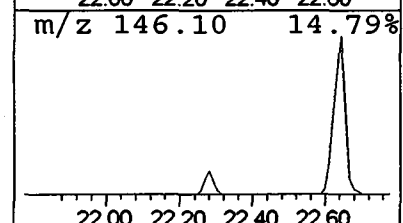
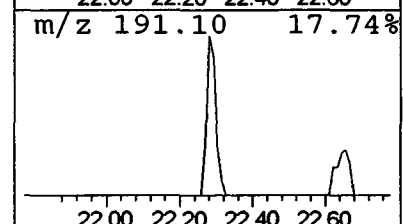
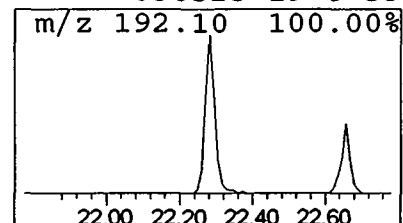
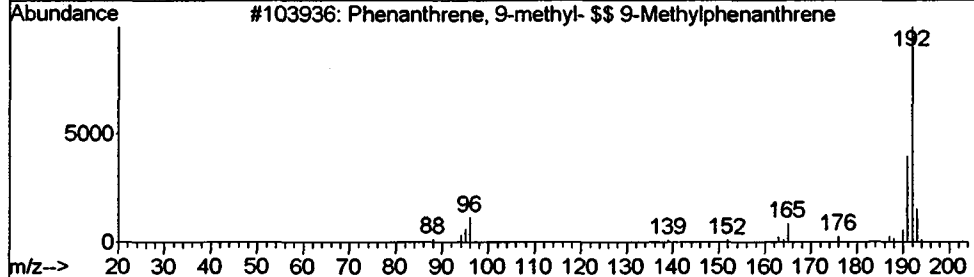
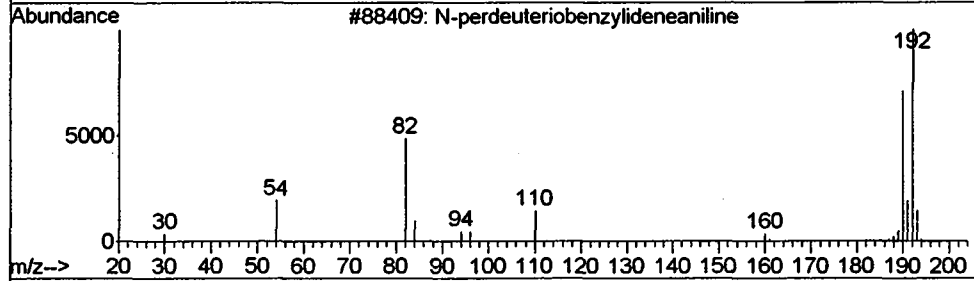
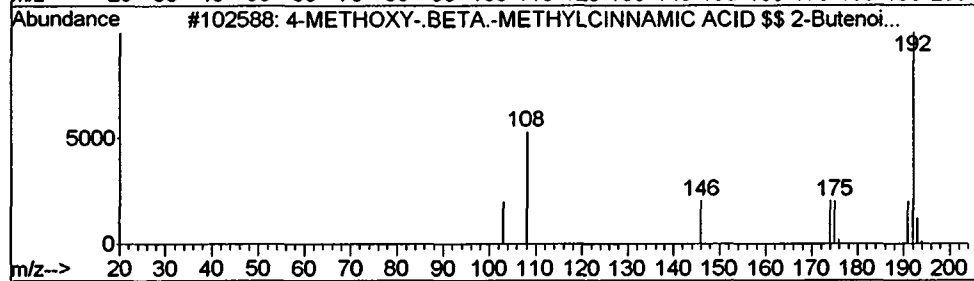
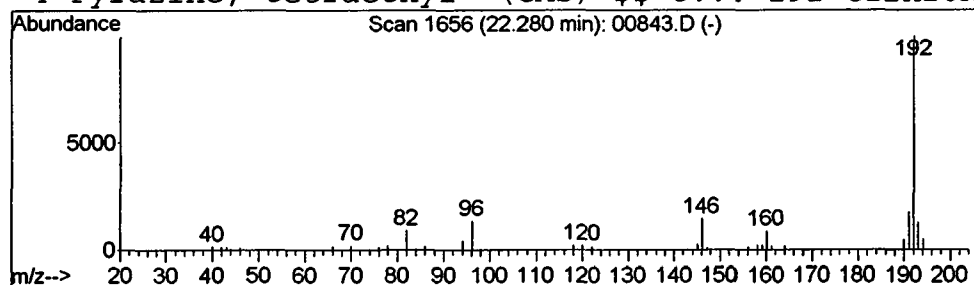
Title : 508 scan method

Library : C:\DATABASE\WILEY7N.L

Peak Number 10 4-METHOXY-.BETA.-METHYLCINN... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.28	0.16 ug/L	150114	Phenanthrene-d10	22.65

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			4-METHOXY-.BETA.-METHYLCINNAMIC ...	192	C11H12O3	019169-94-9	72
2			N-perdeuteriobenzylideneaniline	192	C13D11N	000538-51-2	64
3			Phenanthrene, 9-methyl- \$\$ 9-Met...	192	C15H12	000883-20-5	59
4			Pyrazine, tetraethyl- (CAS) \$\$ t...	192	C12H20N2	038325-19-8	53



Library Search Compound Report

Data File : C:\MSDCHEM\1\5973N\02JAN31\00843.D
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 Sample : 508 scan
 Misc : January 31, 2002
 MS Integration Params: LSCINT.P

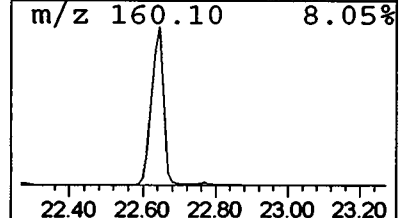
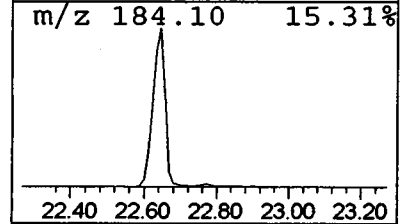
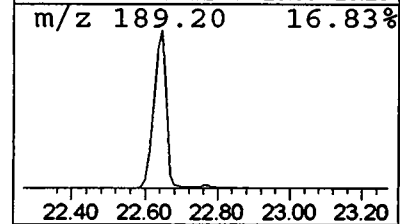
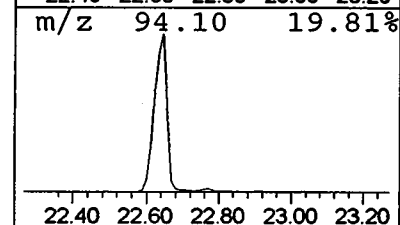
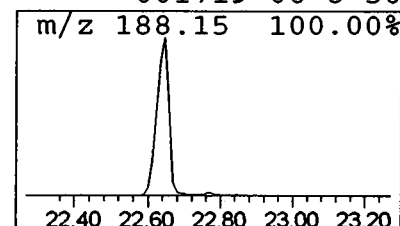
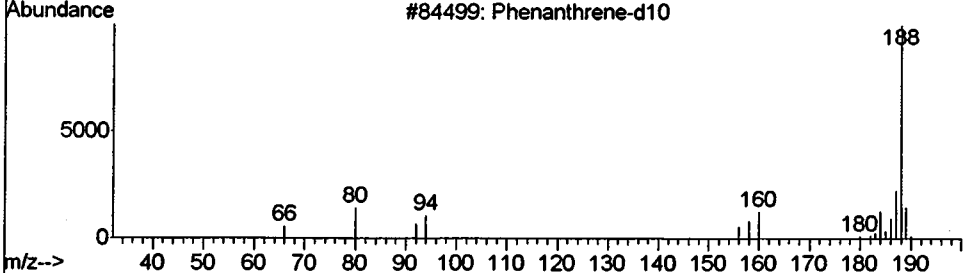
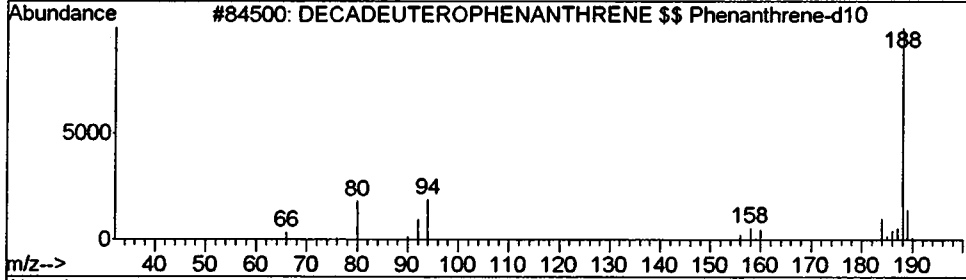
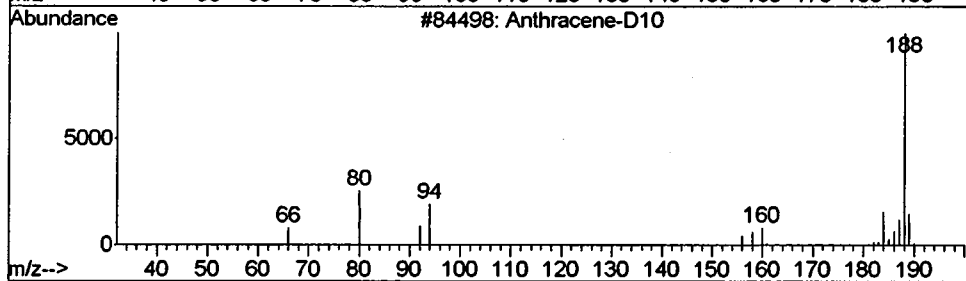
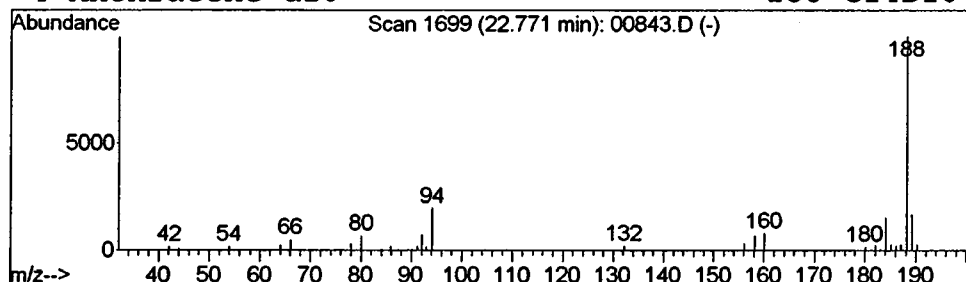
Vial: 5
 Operator:
 Inst : Instrumen
 Multiplr: 1.00

Quant Method : C:\MSDCHEM\1...\SCAN508.M (RTE Integrator)
 Title : 508 scan method
 Library : C:\DATABASE\WILEY7N.L

 Peak Number 11 Anthracene-D10 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.77	0.49 ug/L	443596	Phenanthrene-d10	22.65

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Anthracene-D10	188	C14D10	000000-00-0	90
2	DECADEUTEROPHENANTHRENE \$\$ Phena...	188	C14D10	001517-22-2	86
3	Phenanthrene-d10	188	C14D10	001517-22-2	64
4	Anthracene-d10-	188	C14D10	001719-06-8	50



Library Search Compound Report

Data File : C:\MSDCHEM\1\5973N\02JAN31\00843.D
Acq On : 31 Jan 2002 6:01 pm
Sample : 508 scan
Misc : January 31, 2002
MS Integration Params: LSCINT.P

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Operator:
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Multiplr: 1.00

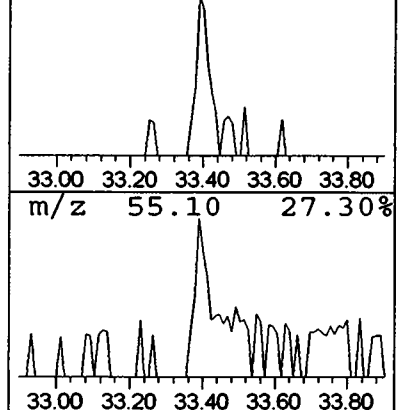
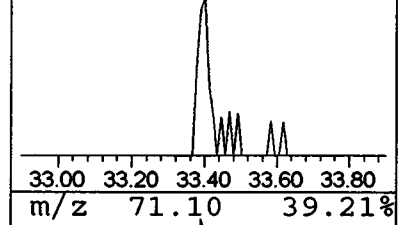
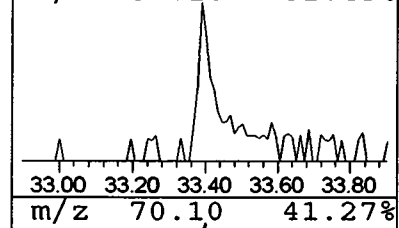
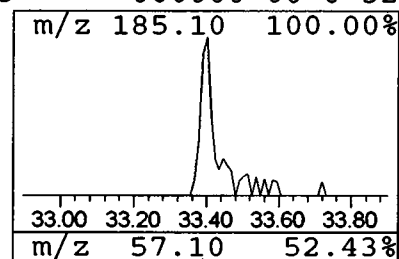
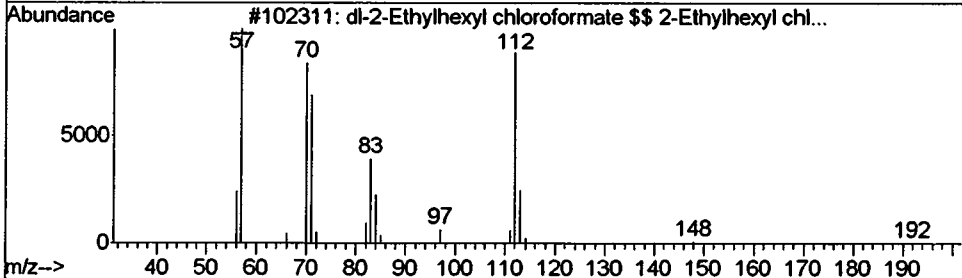
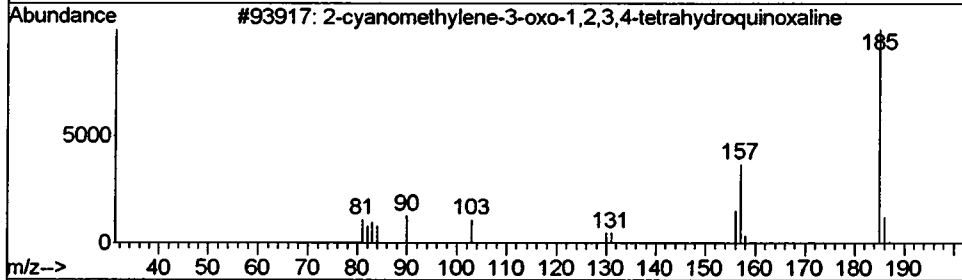
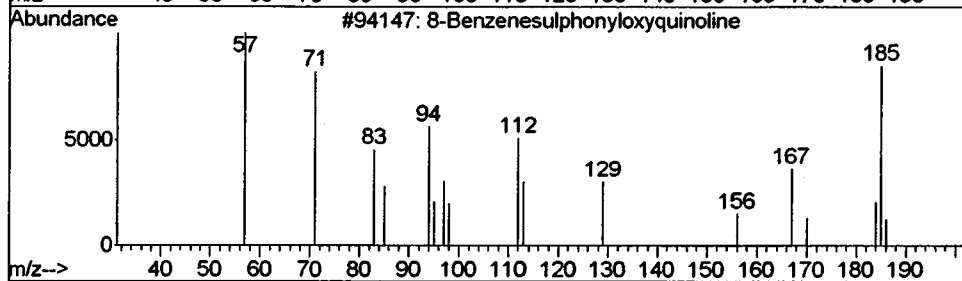
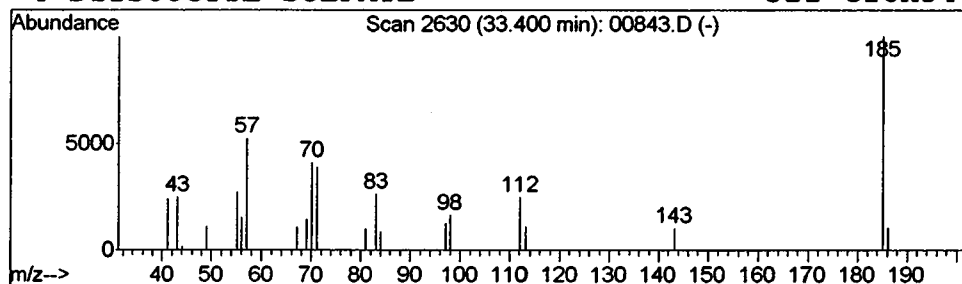
Quant Method : C:\MSDCHEM\1...\SCAN508.M (RTE Integrator)
Title : 508 scan method
Library : C:\DATABASE\WILEY7N.L

Peak Number 12 8-Benzenesulphonyloxyquinoline Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
33.40	0.13 ug/L	94750	Perylene-d12	34.42

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			8-Benzenesulphonyloxyquinoline	185	C12H11NO	000000-00-0	50
2			2-cyanomethylene-3-oxo-1,2,3,4-t...	185	C10H7N3O	000000-00-0	38
3			dl-2-Ethylhexyl chloroformate \$\$...	192	C9H17ClO2	024468-13-1	35
4			DIISOCTYL SULFATE	322	C16H34O4S	000000-00-0	32

NO
GOOD
MATCH



Tentatively Identified Compound (LSC) summary

Operator ID: Date Acquired: 31 Jan 2002 6:01 pm
 Data File: C:\MSDCHEM\1\5973N\02JAN31\00843.D
 Name: 508 scan
 Misc: January 31, 2002
 Method: C:\MSDCHEM\1\METHODS\METHODS\SCAN508.M (RTE Integrator)
 Title: 508 scan method
 Library Searched: C:\DATABASE\WILEY7N.L

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Dichloroacetonitr...	3.50	0.1 ug/L		45160	1	9.71	11147500	20.0
Butanoic acid, me...	3.62	0.1 ug/L		60757	1	9.71	11147500	20.0
2,2-Dimethoxybutane	4.25	0.2 ug/L		84914	1	9.71	11147500	20.0
1-Hexene, 6-bromo...	4.87	0.1 ug/L		42047	1	9.71	11147500	20.0
3-(CYCLOHEX-3'-EN...	5.93	0.3 ug/L		189474	1	9.71	11147500	20.0
2-Propanone, 1,1,...	6.01	0.1 ug/L		71335	1	9.71	11147500	20.0
2-Pentanone, 4-hy...	6.06	0.2 ug/L		91643	1	9.71	11147500	20.0
2-Propanol, 1-(2-...	9.87	0.1 ug/L		47781	1	9.71	11147500	20.0
2,3,4,5-Tetrahydr...	11.27	0.1 ug/L		45385	1	9.71	11147500	20.0
4-METHOXY-.BETA.-...	22.28	0.2 ug/L		150114	4	22.65	18244800	20.0
Anthracene-D10	22.77	0.5 ug/L		443596	4	22.65	18244800	20.0
8-Benzenesulphony...	33.40	0.1 ug/L		94750	6	34.42	14125100	20.0