

User: Vinci, David L
Westchester Dept. of Labs & Research
Date: 01/28/2002 Time: 09:31:32

Sample: AE00168
Analysis: \$GCMS GCMS Scan For Unknowns
Units: ug
Started: 1/28/02 09:31 Ended: 1/28/02 09:31 Analyst: DLV
Page: 1

	Component Name	Viol	Result	Qualifier	MDL
1	Other unknown compounds		.		
2	Surr_2,4-DDD		70		5.0

Comments for Sample AE00168 - Analysis \$GCMS

Westchester Dept. of Labs & Research

User: Vinci, David L

Date: 01-28-2002 Time: 09:31:43

The GCMS scan, from 35 to 550 amu, does not reveal the presence of unusual compounds.

Data File : C:\MSDCHEM\1\5973N\02JAN24\00168.D Vial: 1
 Acq On : 24 Jan 2002 2:59 pm Operator:
 Sample : 508scan Inst : Instrumen
 Misc : January 24, 2002 Multiplr: 1.00
 MS Integration Params: SPLIT.P
 Quant Time: Jan 25 10:22:03 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	9.72	152	1491818	20.00	ug/L	0.00
10) Naphthalene-d8	13.19	136	6431583	20.00	ug/L	-0.02
17) Acenaphthene-d10	18.34	164	3332964	20.00	ug/L	-0.01
29) Phenanthrene-d10	22.65	188	5447319	20.00	ug/L	-0.02
38) Chrysene-d12	30.50	240	4440597	20.00	ug/L	-0.01
44) Perylene-d12	34.42	264	3175060	20.00	ug/L	0.00

System Monitoring Compounds		R.T.	QIon	Response	Conc	Units	Dev (Min)
37) 2,4-DDD		27.73	235	301692	3.50	ug/L	0.00
Spiked Amount	5.000			Recovery	=	70.00%	

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
20) Dimethylphthalate	18.34	163	533948	2.43	ug/L	# 56
21) 2,6-Dinitrotoluene	18.34	165	416775	9.71	ug/L	# 17
42) Bis(2-ethylhexyl)phthalate	31.11	149	10370	0.60	ug/L	81

 (#) = qualifier out of range (m) = manual integration
 00168.D SCAN508.M Fri Jan 25 10:22:04 2002

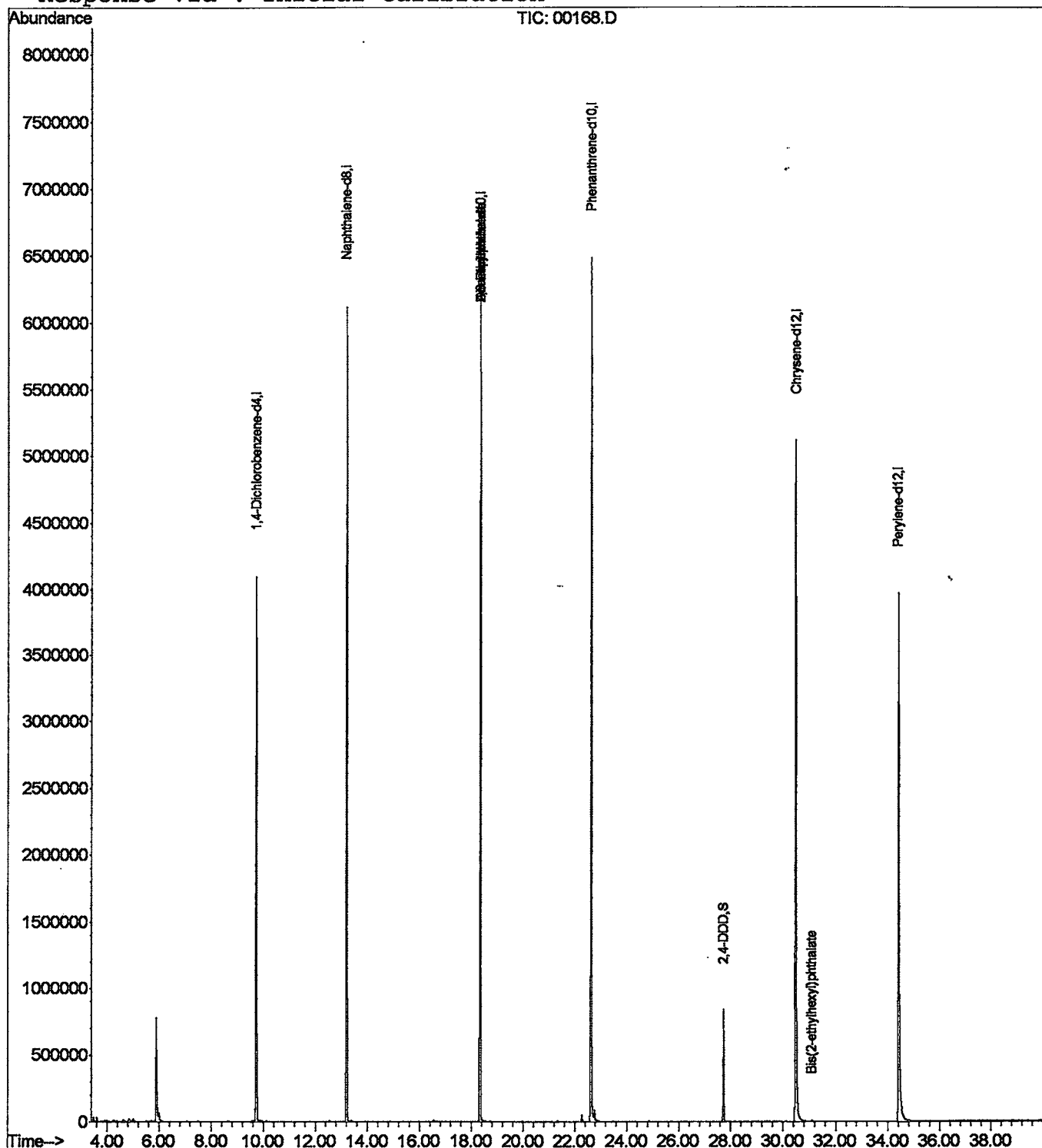
Page 1

Data File : C:\MSDCHEM\1\5973N\02JAN24\00168.D
 Acq On : 24 Jan 2002 2:59 pm
 Sample : 508scan
 Misc : January 24, 2002
 MS Integration Params: SPLIT.P
 Quant Time: Jan 25 10:22 2002

Vial: 1
 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\02JAN24\00168.D

Vial: 1

Acq Ori : 24 Jan 2002 2:59 pm

Operator:

Sample : 508scan

Inst : Instrumen

Misc : January 24, 2002

Multiplr: 1.00

MS Integration Params: LSCINT.P

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

Title : 508 scan method

Smoothing : OFF

Filtering: 5

Sampling : 2

Min Area: 100 Area counts

Start Thrs: 0.2

Max Peaks: 20

Stop Thrs : 0

Peak Location: TOP

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

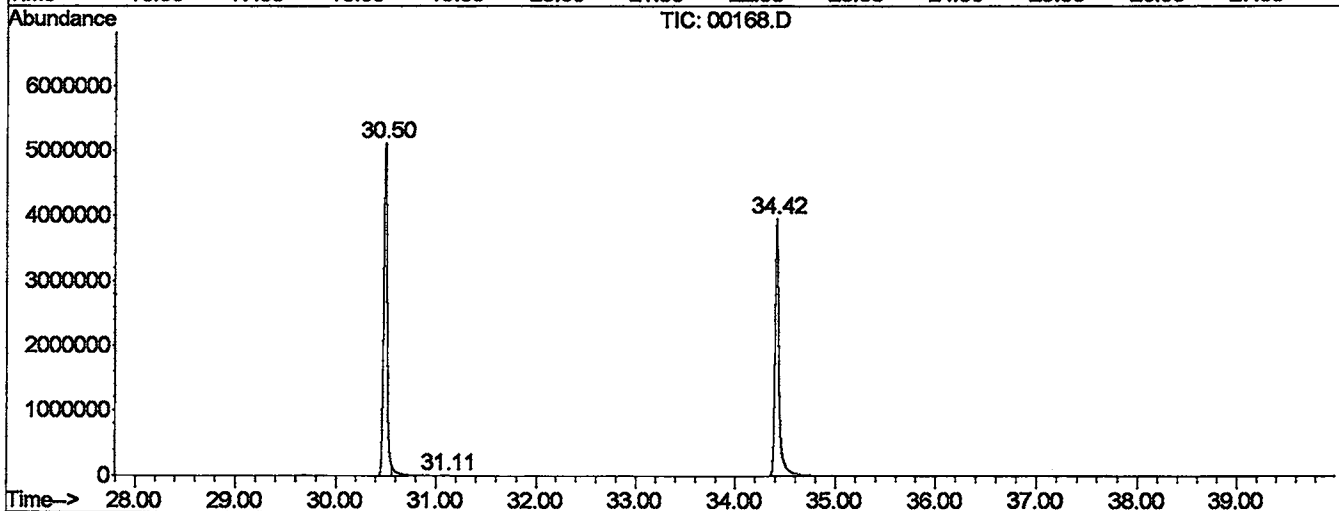
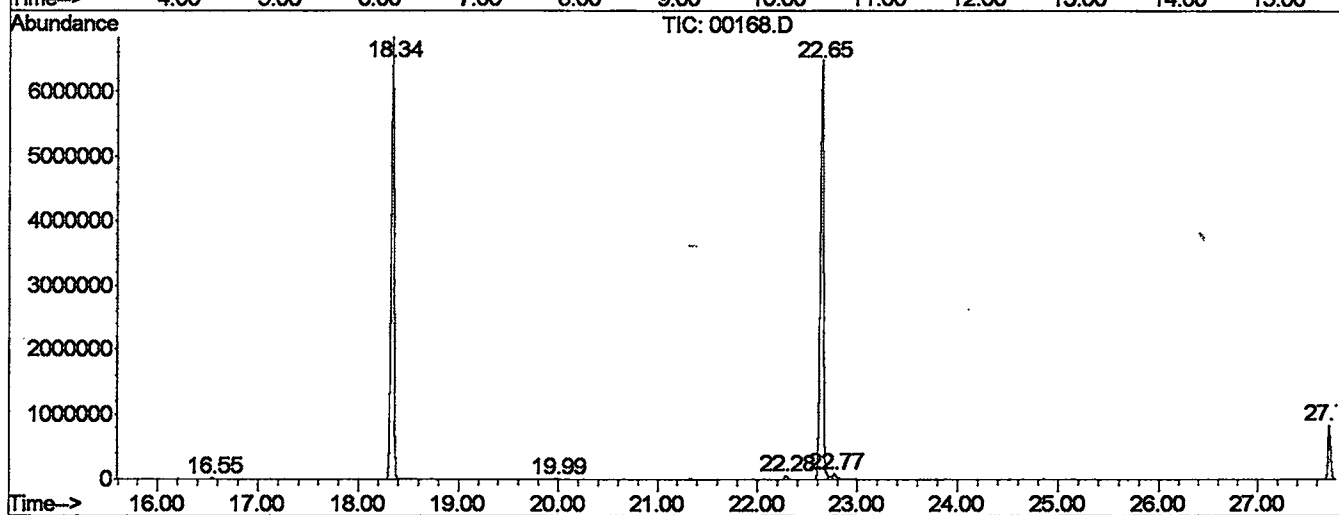
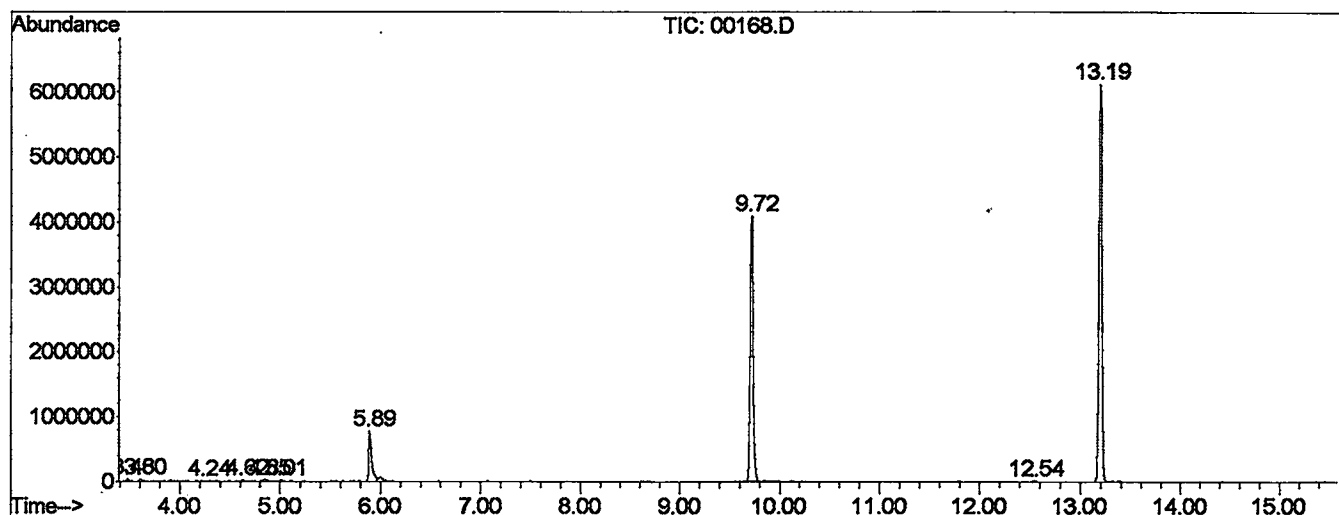
Signal : TIC

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.476	7	9	13	rVB	31653	43334	0.30%	0.057%
2	3.602	15	20	25	rBV	33853	48845	0.34%	0.064%
3	4.241	73	76	85	rVB5	9645	34250	0.24%	0.045%
4	4.618	107	109	115	rBV2	16331	40152	0.28%	0.053%
5	4.846	123	129	141	rBV4	25036	95892	0.66%	0.126%
6	5.006	141	143	149	rVB	23249	45260	0.31%	0.059%
7	5.885	217	220	227	rBV	776524	1597408	11.02%	2.093%
8	9.721	551	556	563	rVV	4093382	8505727	58.70%	11.142%
9	12.541	799	803	809	rVB	13952	23493	0.16%	0.031%
10	13.192	855	860	865	rBV	6121921	12392333	85.52%	16.234%
11	16.549	1149	1154	1159	rVV2	16362	30979	0.21%	0.041%
12	18.341	1305	1311	1315	rBV	6833629	13563239	93.60%	17.768%
13	19.985	1445	1455	1459	rVB2	7400	31235	0.22%	0.041%
14	22.280	1651	1656	1661	rVB	54823	101559	0.70%	0.133%
15	22.645	1681	1688	1693	rBV2	6485545	14490042	100.00%	18.982%
16	22.771	1697	1699	1719	rVB	83268	234038	1.62%	0.307%
17	27.726	2129	2133	2139	rVV	835729	1491988	10.30%	1.954%
18	30.500	2369	2376	2381	rBV	5117968	12714680	87.75%	16.656%
19	31.105	2423	2429	2437	rBV4	9075	28980	0.20%	0.038%
20	34.416	2713	2719	2753	rBV	3964662	10822820	74.69%	14.178%

Sum of corrected areas: 76336254

LSC Report - Integrated Chromatogram

File : C:\MSDCHEM\1\5973N\02JAN24\00168.D
 Operator :
 Acquired : 24 Jan 2002 2:59 pm using AcqMethod SCAN508
 Instrument : Instrumen
 Sample Name: 508scan
 Misc Info : January 24, 2002
 Vial Number: 1
 Quant File :SCAN508.RES (RTE Integrator)



Library Search Compound Report

Data File : C:\MSDCHEM\1\5973N\02JAN24\00168.D
 Acq On : 24 Jan 2002 2:59 pm
 Sample : 508scan
 Misc : January 24, 2002
 MS Integration Params: LSCINT.P

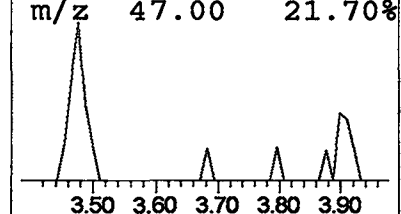
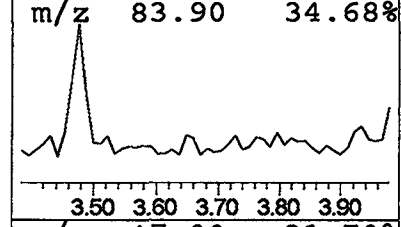
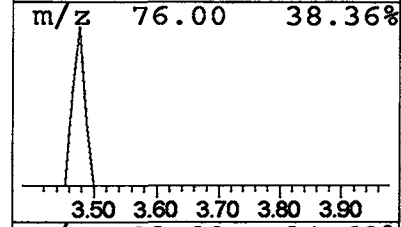
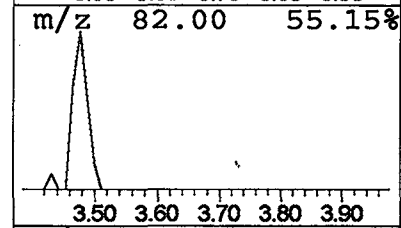
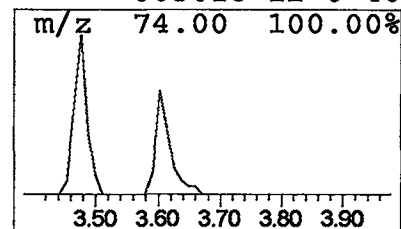
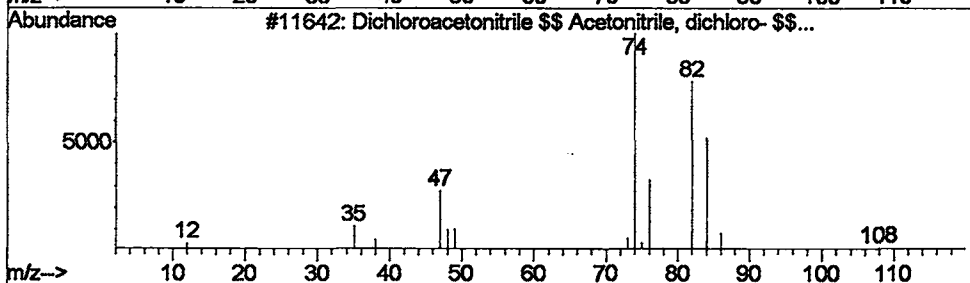
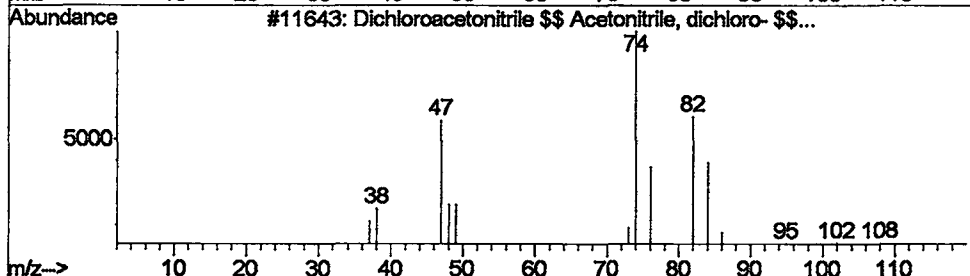
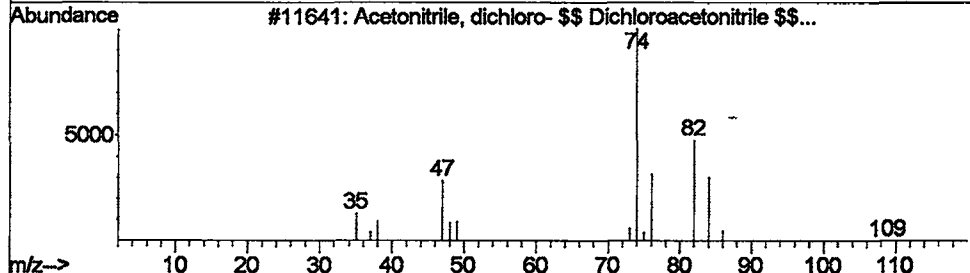
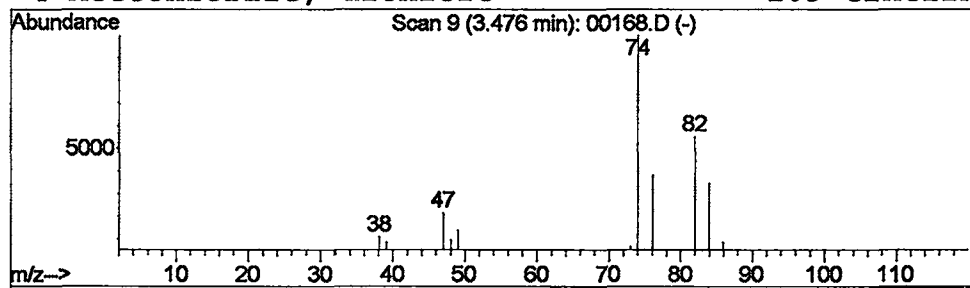
Vial: 1
 Operator:
 Inst : Instrumen
 Multiplr: 1.00

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

 Peak Number 1 Acetonitrile, dichloro- \$\$... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.48	0.10 ug/L	43334	1,4-Dichlorobenzene-d4	9.72

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Acetonitrile, dichloro- \$\$ Dichl...	109	C2HCl2N	003018-12-0	83
2		Dichloroacetonitrile \$\$ Acetonit...	109	C2HCl2N	003018-12-0	83
3		Dichloroacetonitrile \$\$ Acetonit...	109	C2HCl2N	003018-12-0	53
4		Acetonitrile, dichloro-	109	C2HCl2N	003018-12-0	40



Library Search Compound Report

Data File : C:\MSDCHEM\1\5973N\02JAN24\00168.D
Acq On : 24 Jan 2002 2:59 pm
Sample : 508scan
Misc : January 24, 2002
MS Integration Params: LSCINT.P

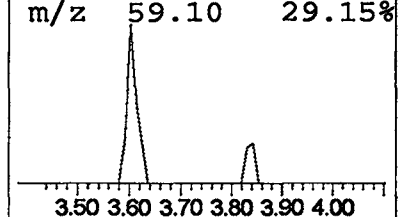
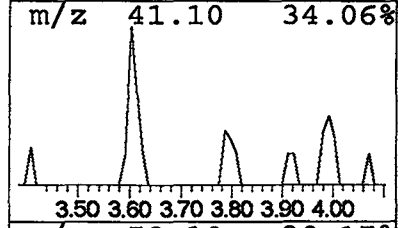
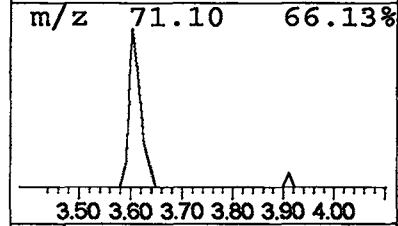
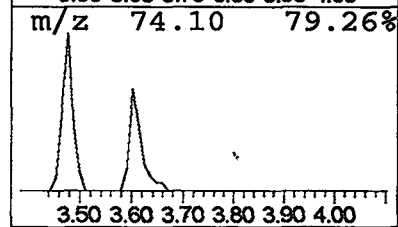
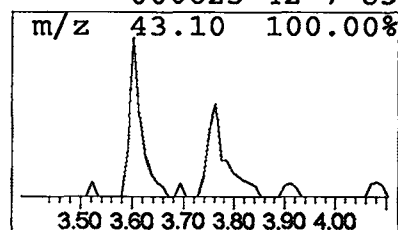
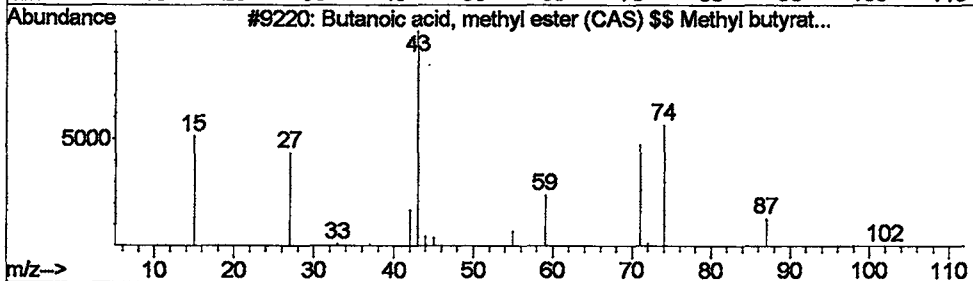
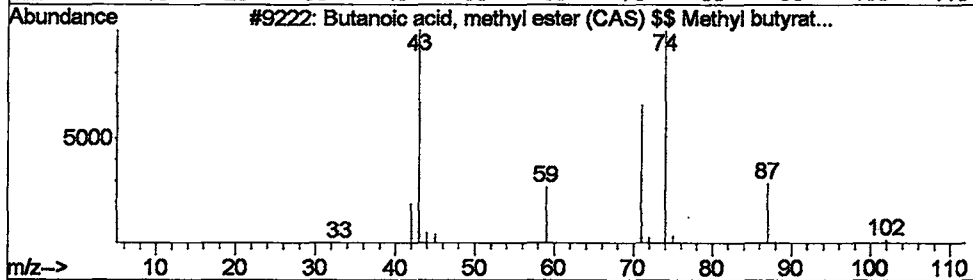
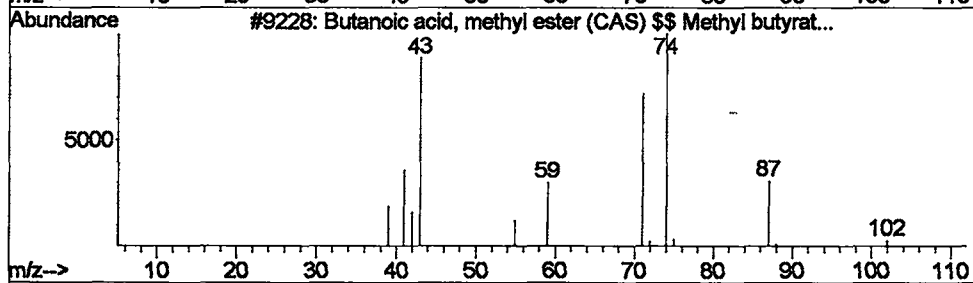
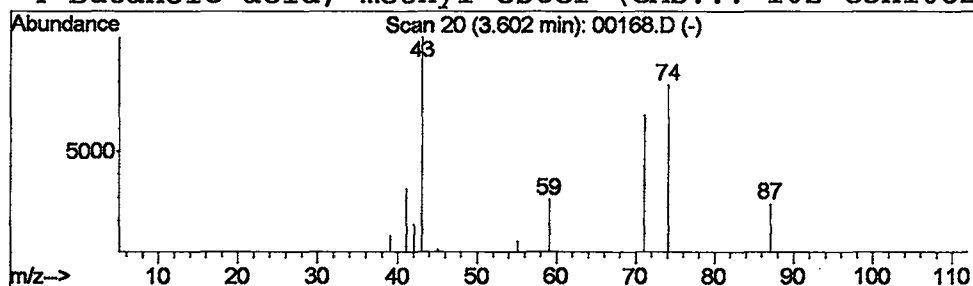
Vial: 1
Operator:
Inst : Instrumen
Multiplr: 1.00

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 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.60	0.11 ug/L	48845	1,4-Dichlorobenzene-d4	9.72

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Butanoic acid, methyl ester (CAS...	102	C5H10O2	000623-42-7	83
2		Butanoic acid, methyl ester (CAS...	102	C5H10O2	000623-42-7	83
3		Butanoic acid, methyl ester (CAS...	102	C5H10O2	000623-42-7	83
4		Butanoic acid, methyl ester (CAS...	102	C5H10O2	000623-42-7	83



Library Search Compound Report

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Acq On : 24 Jan 2002 2:59 pm

Sample : 508scan

Misc : January 24, 2002

MS Integration Params: LSCINT.P

Vial: 1

Operator:

Inst : Instrumen

Multiplr: 1.00

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

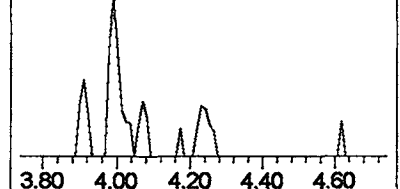
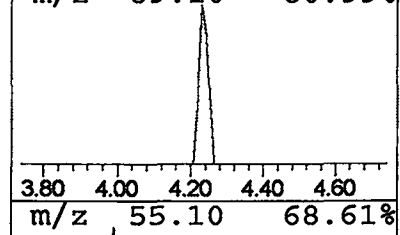
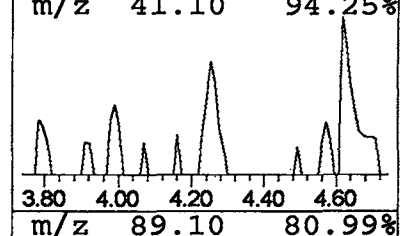
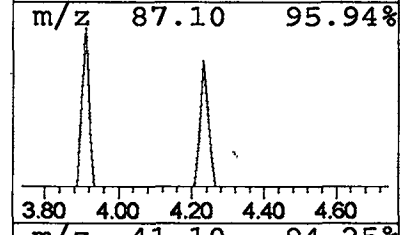
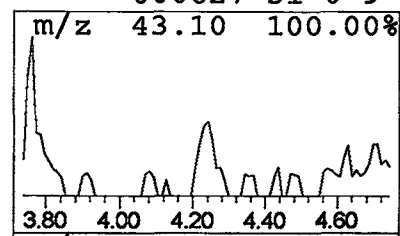
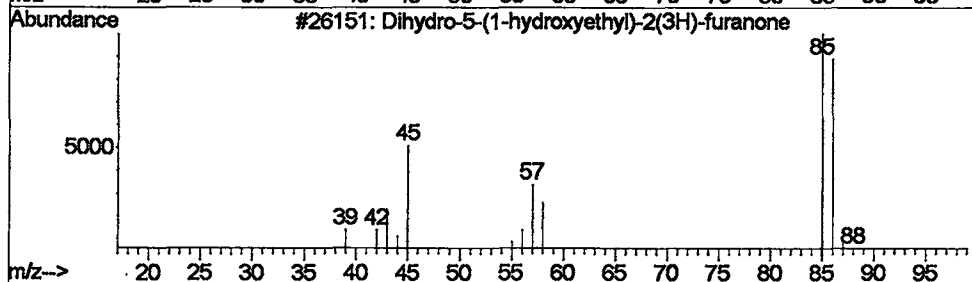
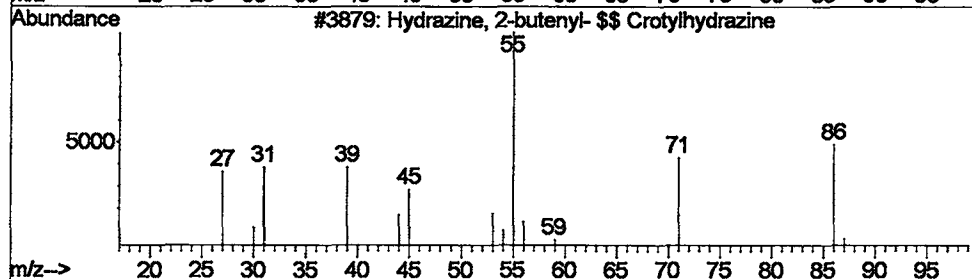
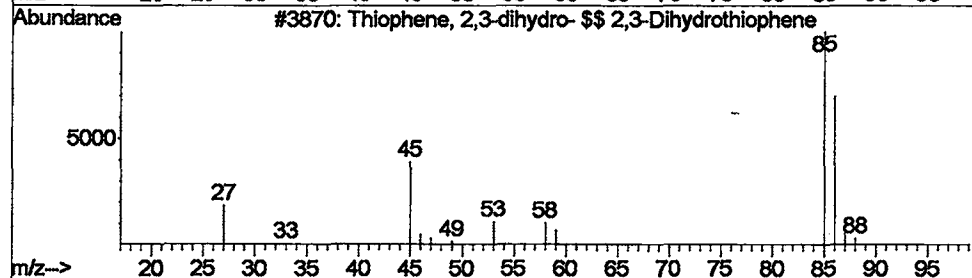
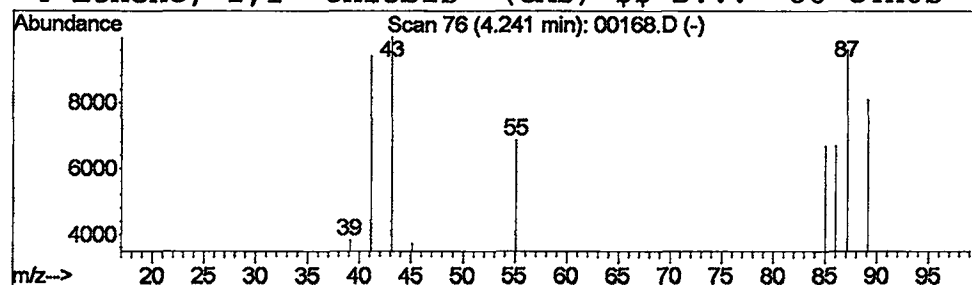
Title : 508 scan method

Library : C:\DATABASE\WILEY7N.L

Peak Number 3 Thiophene, 2,3-dihydro- \$\$... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.24	0.08 ug/L	34250	1,4-Dichlorobenzene-d4	9.72

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Thiophene, 2,3-dihydro- \$\$ 2,3-D...	86	C4H6S	001120-59-8	9
2		Hydrazine, 2-butenyl- \$\$ Crotylh...	86	C4H10N2	036566-68-4	9
3		Dihydro-5-(1-hydroxyethyl)-2(3H)... 130	130	C6H10O3	000000-00-0	9
4		Ethene, 1,1'-thiobis- (CAS) \$\$ D...	86	C4H6S	000627-51-0	9



Library Search Compound Report

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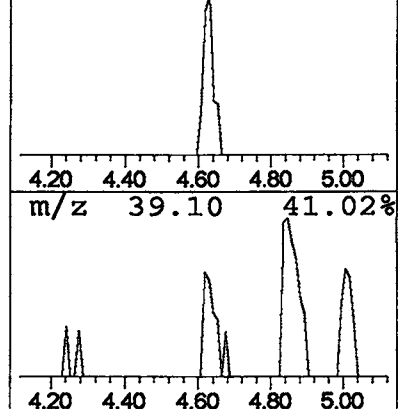
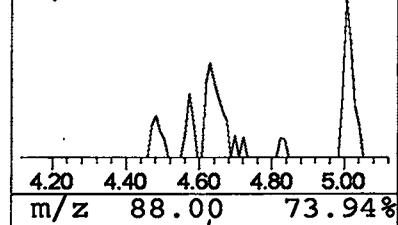
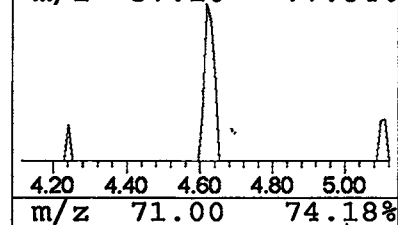
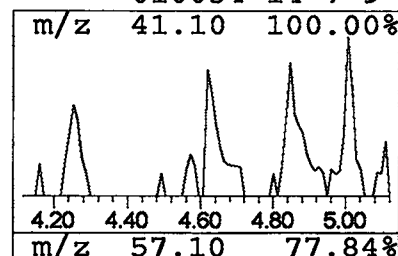
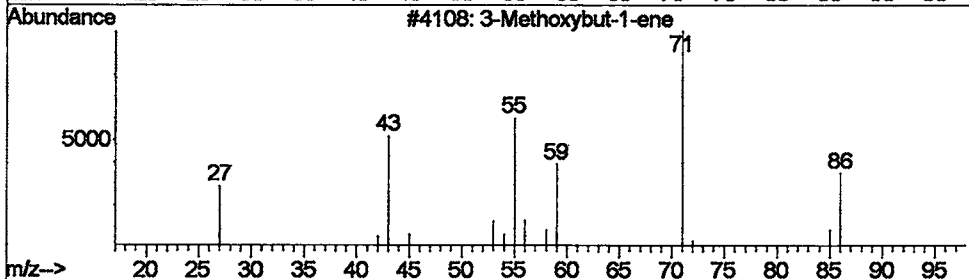
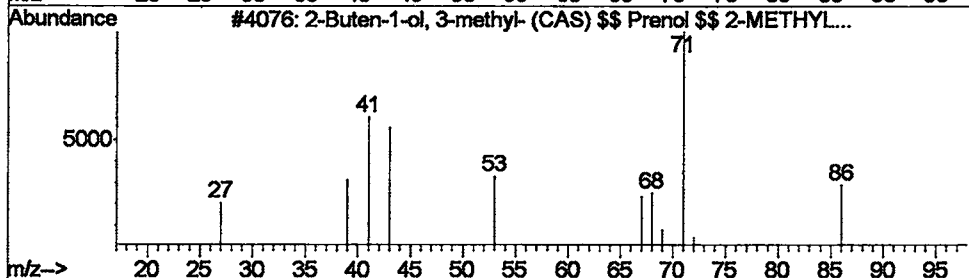
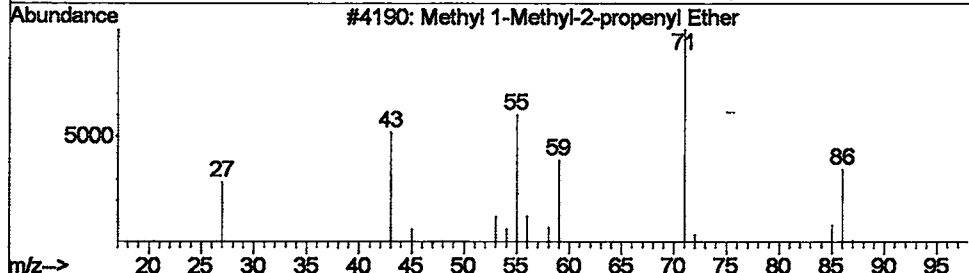
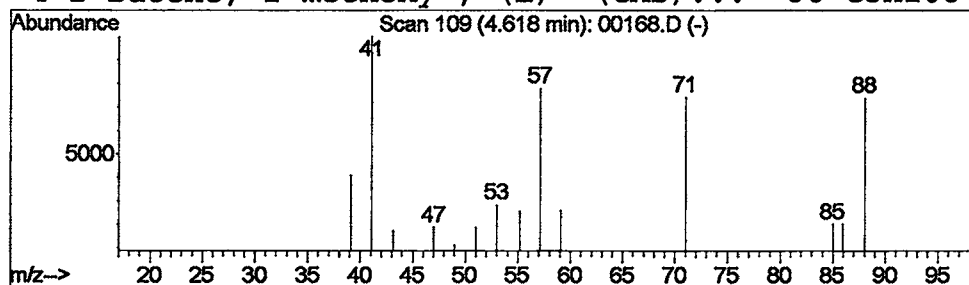
Vial: 1
 Operator:
 Inst : Instrumen
 Multiplr: 1.00

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

 Peak Number 4 Methyl 1-Methyl-2-propenyl ... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.62	0.09 ug/L	40152	1,4-Dichlorobenzene-d4	9.72

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Methyl 1-Methyl-2-propenyl Ether	86	C5H10O	000000-00-0	9
2			2-Buten-1-ol, 3-methyl- (CAS) \$\$...	86	C5H10O	000556-82-1	9
3			3-Methoxybut-1-ene	86	C5H10O	017351-24-5	9
4			2-Butene, 1-methoxy-, (E)- (CAS)...	86	C5H10O	010034-14-7	9



Library Search Compound Report

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Acq On : 24 Jan 2002 2:59 pm

Sample : 508scan

Misc : January 24, 2002

MS Integration Params: LSCINT.P

Vial: 1

Operator:

Inst : Instrumen

Multiplr: 1.00

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

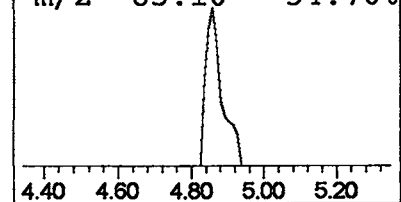
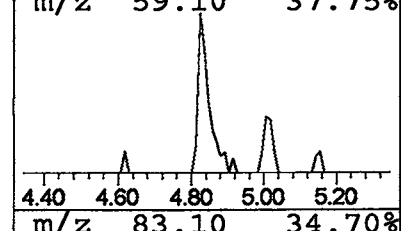
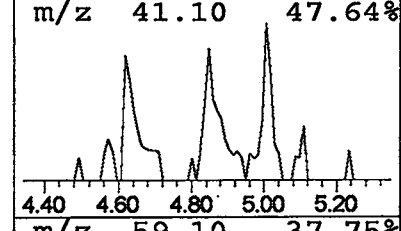
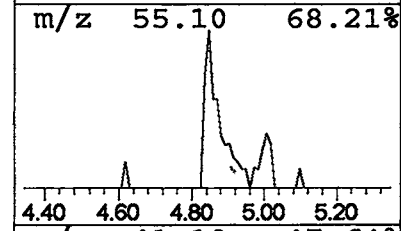
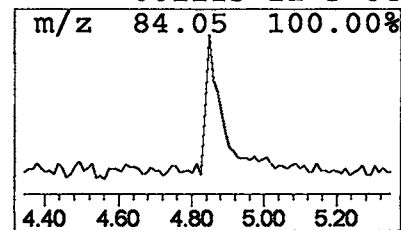
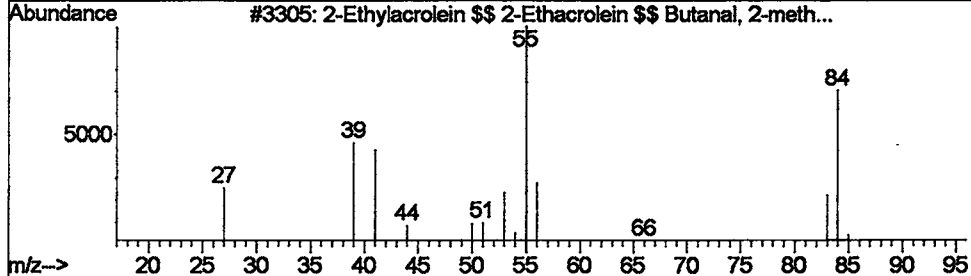
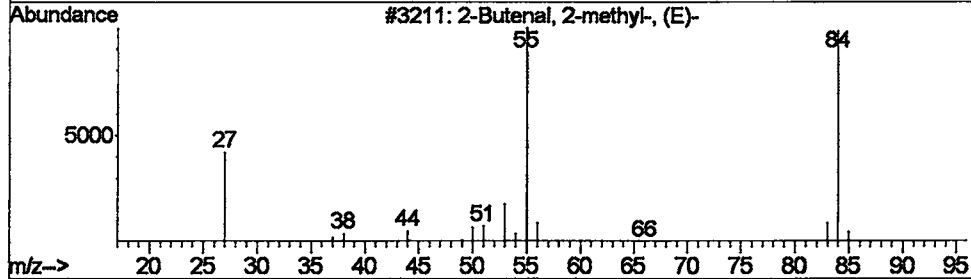
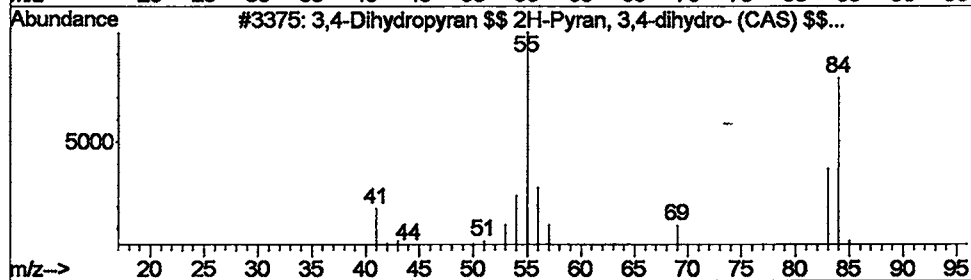
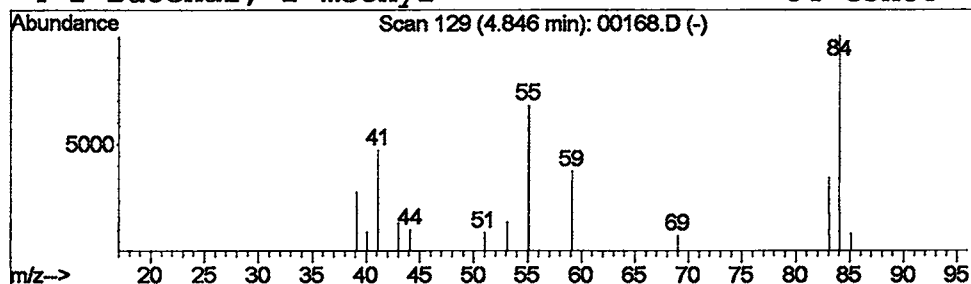
Title : 508 scan method

Library : C:\DATABASE\WILEY7N.L

Peak Number 5 3,4-Dihydropyran \$\$ 2H-Pyra... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.85	0.23 ug/L	95892	1,4-Dichlorobenzene-d4	9.72

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			3,4-Dihydropyran \$\$ 2H-Pyran, 3,...	84	C5H8O	000110-87-2	72
2			2-Butenal, 2-methyl-, (E)-	84	C5H8O	000497-03-0	64
3			2-Ethylacrolein \$\$ 2-Ethacrolein...	84	C5H8O	000922-63-4	64
4			2-Butenal, 2-methyl-	84	C5H8O	001115-11-3	64



Library Search Compound Report

Data File : C:\MSDCHEM\1\5973N\02JAN24\00168.D
 Acq Ori : 24 Jan 2002 2:59 pm
 Sample : 508scan
 Misc : January 24, 2002
 MS Integration Params: LSCINT.P

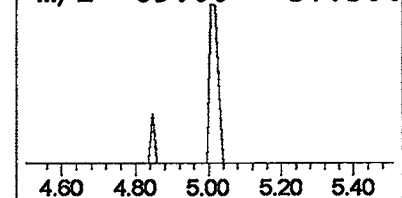
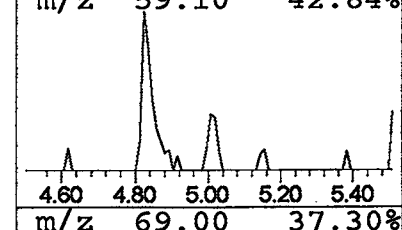
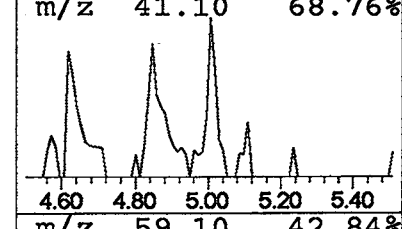
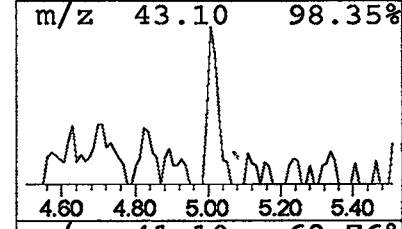
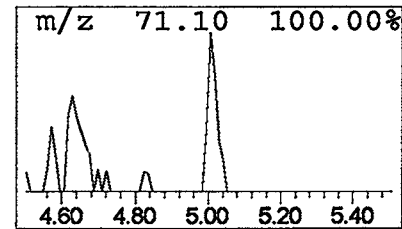
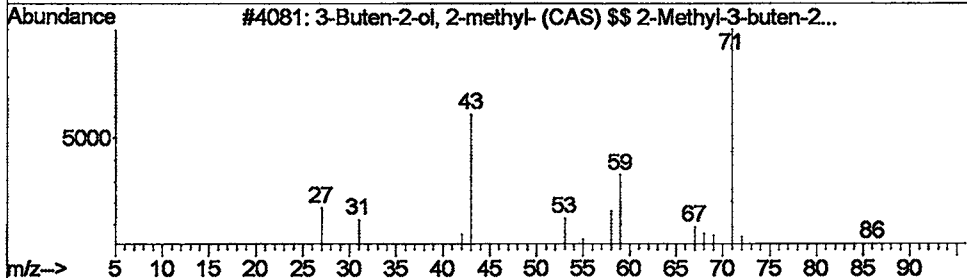
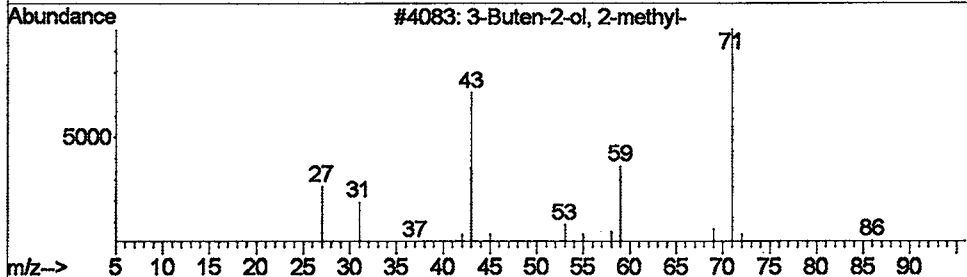
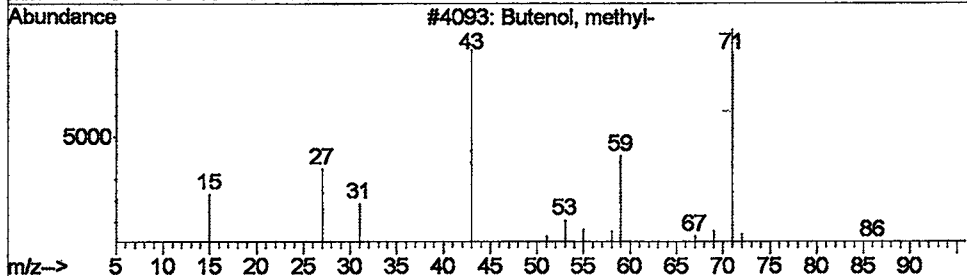
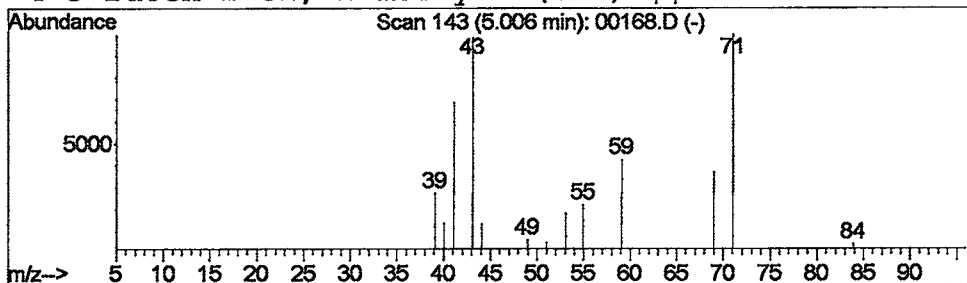
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 Operator:
 Inst : Instrumen
 Multiplr: 1.00

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

 Peak Number 6 Butenol, methyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.01	0.11 ug/L	45260	1,4-Dichlorobenzene-d4	9.72

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Butenol, methyl-	86	C5H10O	060766-00-9	72
2		3-Buten-2-ol, 2-methyl-	86	C5H10O	000115-18-4	56
3		3-Buten-2-ol, 2-methyl- (CAS) \$\$...	86	C5H10O	000115-18-4	50
4		3-Buten-2-ol, 2-methyl- (CAS) \$\$...	86	C5H10O	000115-18-4	9



Library Search Compound Report

Data File : C:\MSDCHEM\1\5973N\02JAN24\00168.D

Acq On : 24 Jan 2002 2:59 pm

Sample : 508scan

Misc : January 24, 2002

MS Integration Params: LSCINT.P

Vial: 1

Operator:

Inst : Instrumen

Multiplr: 1.00

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

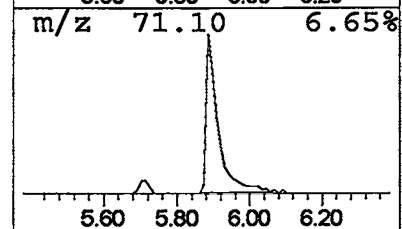
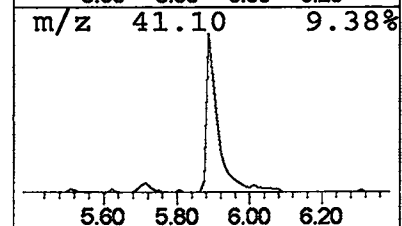
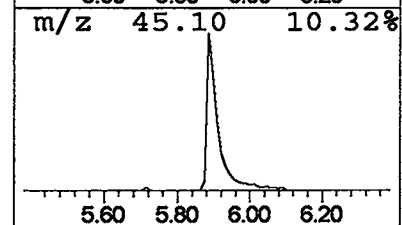
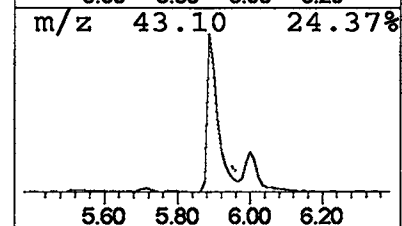
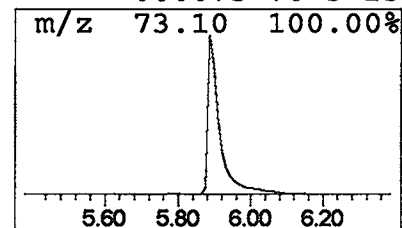
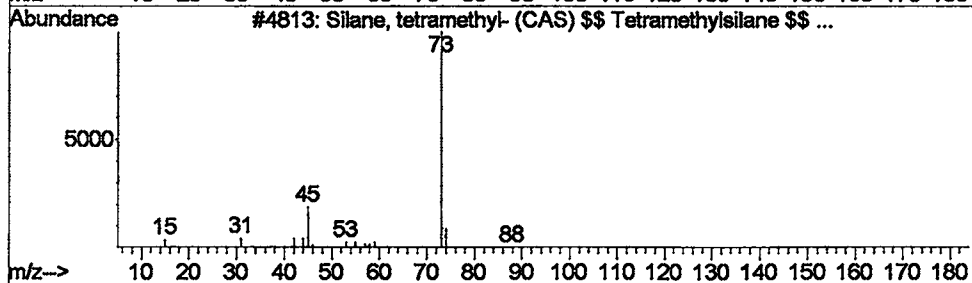
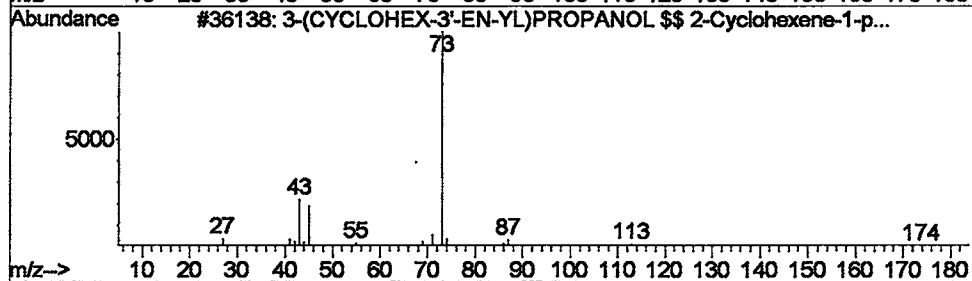
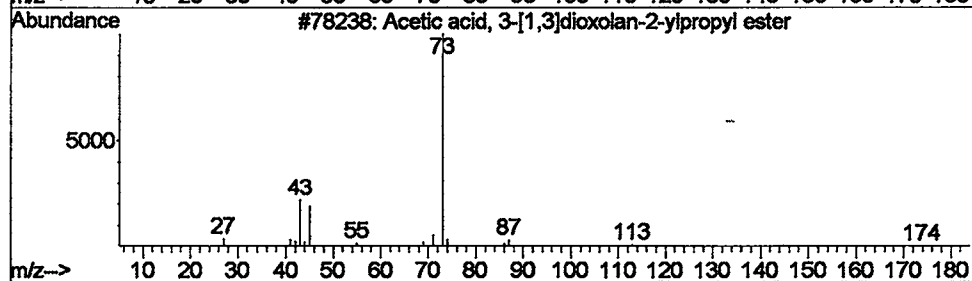
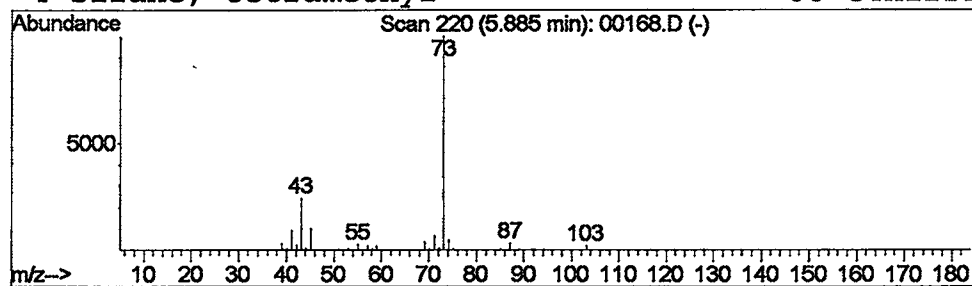
Title : 508 scan method

Library : C:\DATABASE\WILEY7N.L

Peak Number 7 Acetic acid, 3-[1,3]dioxola... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.89	3.76 ug/L	1597410	1,4-Dichlorobenzene-d4	9.72

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Acetic acid, 3-[1,3]dioxolan-2-y...	174	C8H14O4	000000-00-0	33
2			3-(CYCLOHEX-3'-EN-YL)PROPANOL \$\$...	174	C8H14O4	015745-87-6	33
3			Silane, tetramethyl- (CAS) \$\$ Te...	88	C4H12Si	000075-76-3	25
4			Silane, tetramethyl-	88	C4H12Si	000075-76-3	23



Library Search Compound Report

Data File : C:\MSDCHEM\1\5973N\02JAN24\00168.D

Acq On : 24 Jan 2002 2:59 pm

Sample : 508scan

Misc : January 24, 2002

MS Integration Params: LSCINT.P

Vial: 1

Operator:

Inst : Instrumen

Multiplr: 1.00

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

Title : 508 scan method

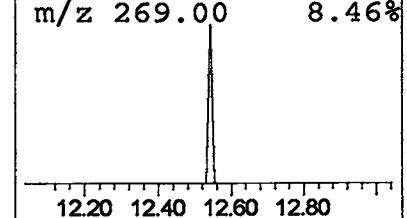
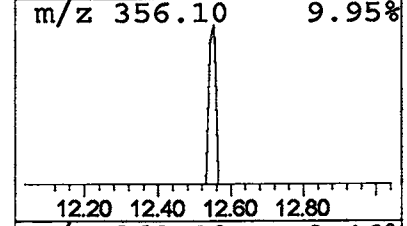
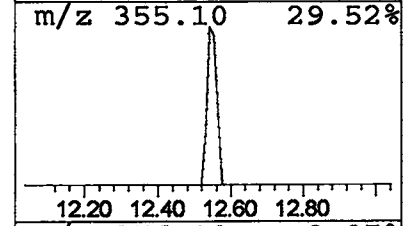
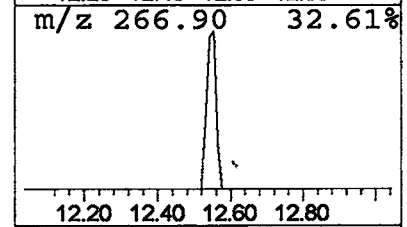
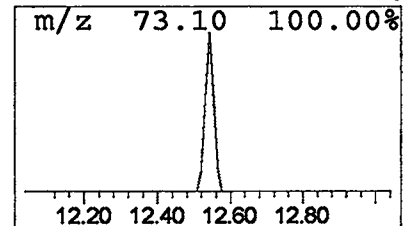
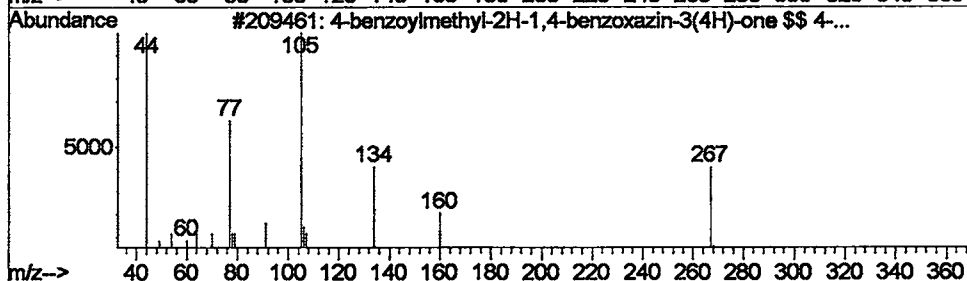
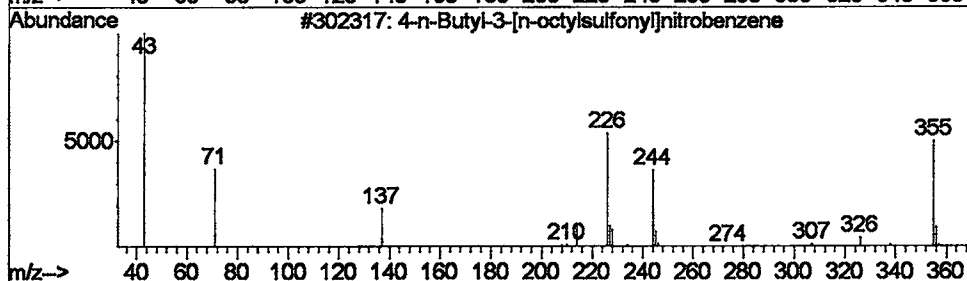
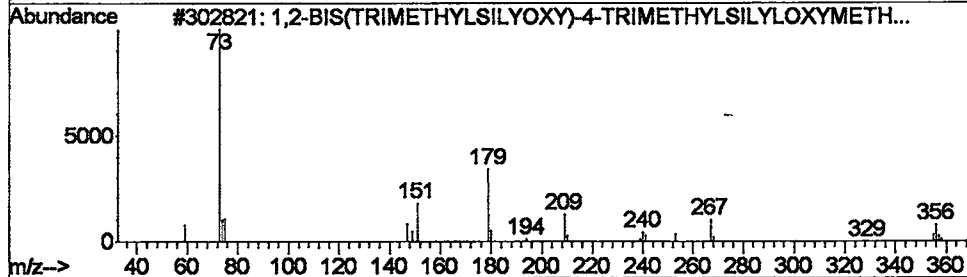
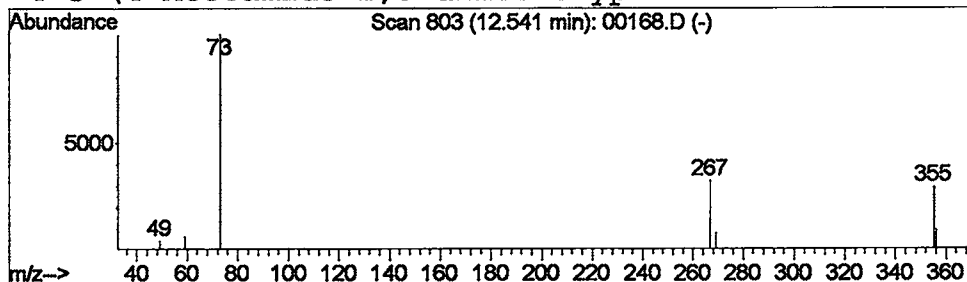
Library : C:\DATABASE\WILEY7N.L

Peak Number 8 1,2-BIS(TRIMETHYLSILYOXY)-4... Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.54	0.04 ug/L	23493	Naphthalene-d8	13.19

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1,2-BIS(TRIMETHYLSILYOXY)-4-TRIM...	356	C16H32O3Si3	000000-00-0	5
2			4-n-Butyl-3-[n-octylsulfonyl]nit...	355	C18H29NO4S	000000-00-0	4
3			4-benzoylmethyl-2H-1,4-benzoxazi...	267	C16H13NO3	105492-43-1	4
4			5-(4-Acetamido-2,3-dimethoxyphen...	355	C20H21NO5	000000-00-0	4

NO
Good
match



Library Search Compound Report

Data File : C:\MSDCHEM\1\5973N\02JAN24\00168.D
Acq On : 24 Jan 2002 2:59 pm
Sample : 508scan
Misc : January 24, 2002
MS Integration Params: LSCINT.P

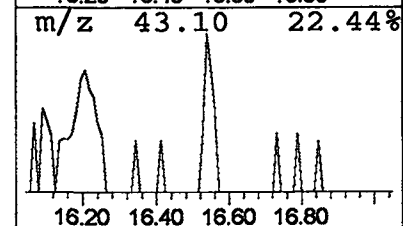
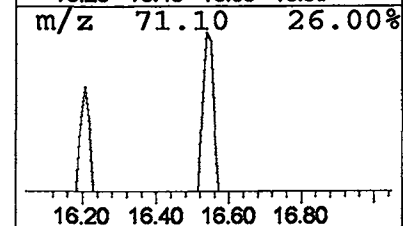
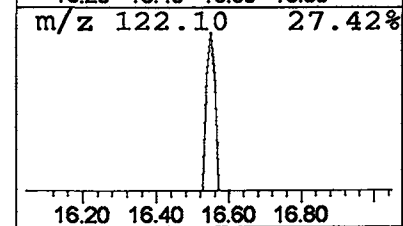
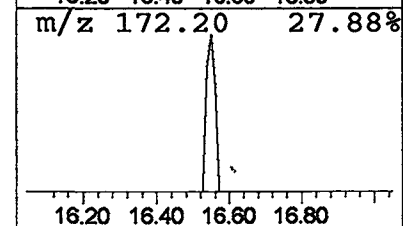
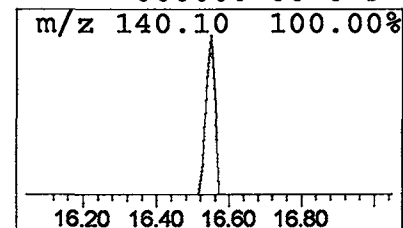
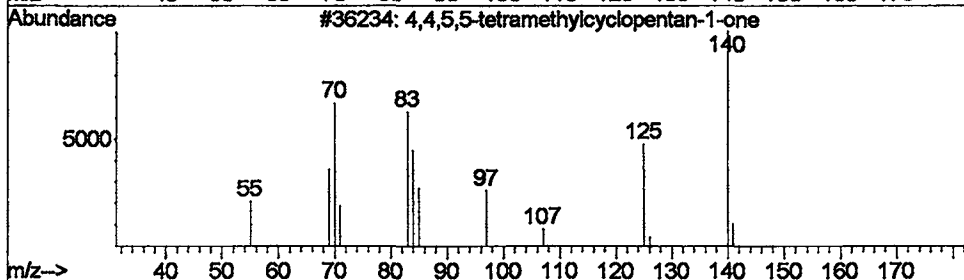
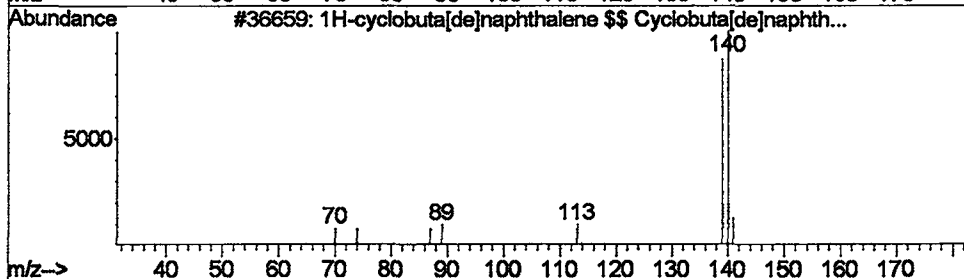
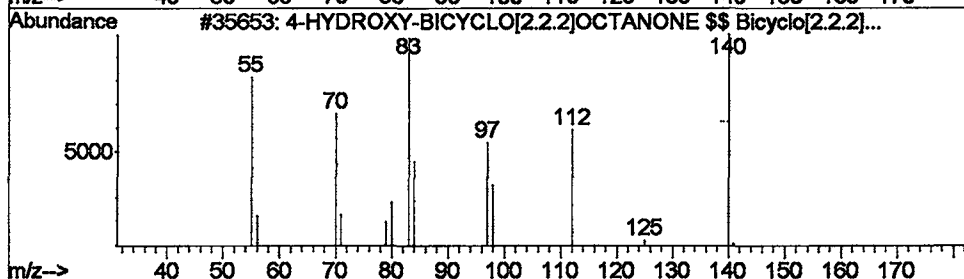
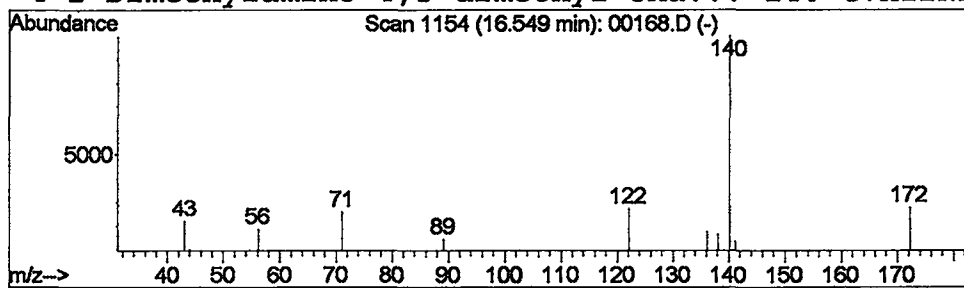
Vial: 1
Operator:
Inst : Instrumen
Multiplr: 1.00

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

Peak Number 9 4-HYDROXY-BICYCLO[2.2.2]OCT... Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.55	0.05 ug/L	30979	Acenaphthene-d10	18.34

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			4-HYDROXY-BICYCLO[2.2.2]OCTANONE...	140	C8H12O2	066348-59-2	38
2			1H-cyclobuta[de]naphthalene \$\$ C...	140	C11H8	024973-91-9	9
3			4,4,5,5-tetramethylcyclopentan-1...	140	C9H16O	111689-30-6	9
4			2-Dimethylamino-4,5-dimethyl-oxa...	140	C7H12N2O	000000-00-0	9



Library Search Compound Report

Data File : C:\MSDCHEM\1\5973N\02JAN24\00168.D
Acq On : 24 Jan 2002 2:59 pm
Sample : 508scan
Misc : January 24, 2002
MS Integration Params: LSCINT.P

Vial: 1
Operator:
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Multiplr: 1.00

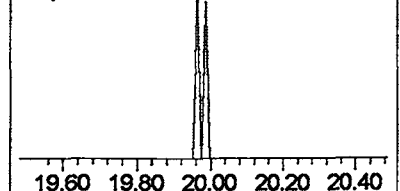
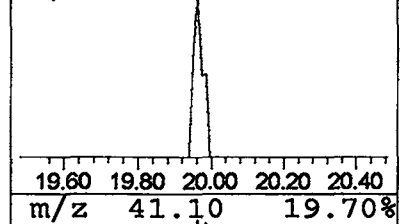
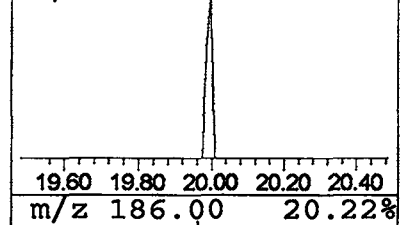
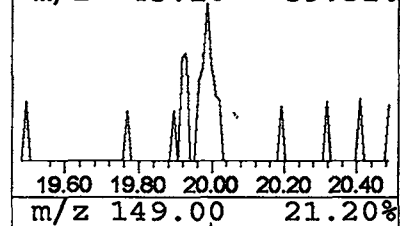
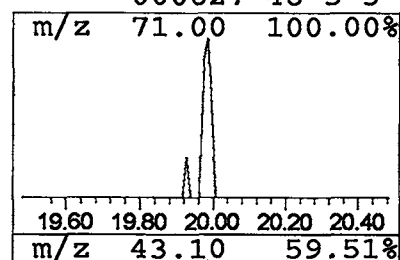
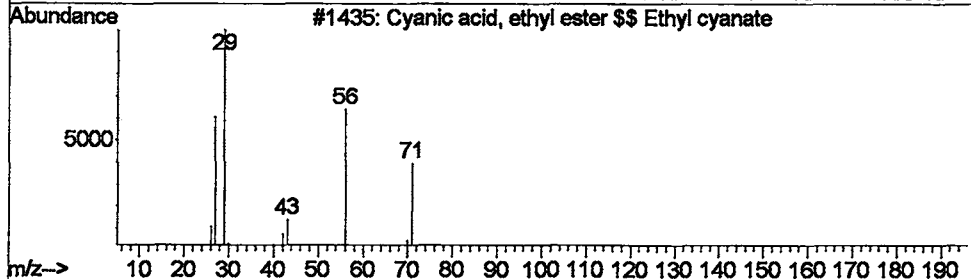
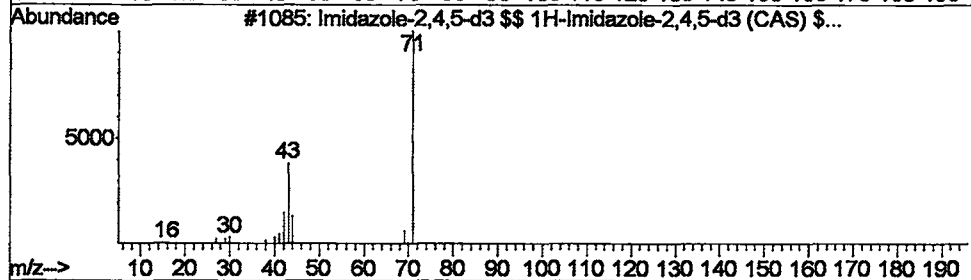
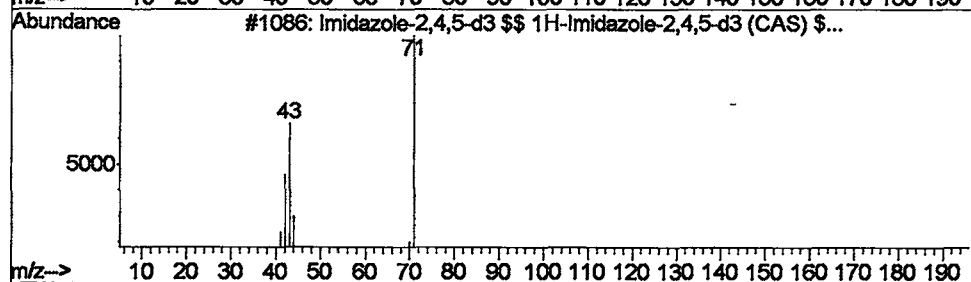
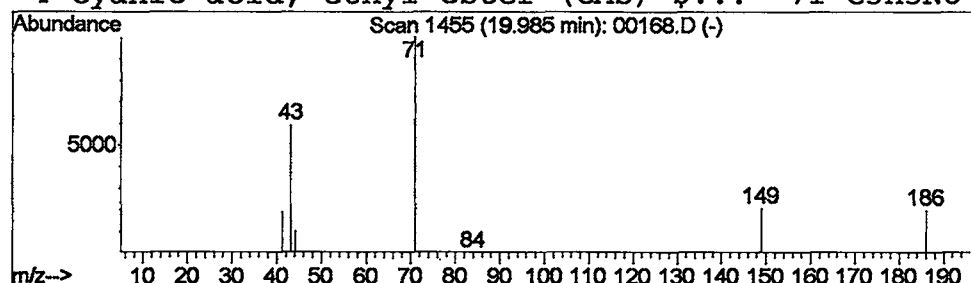
Quant Method : C:\MSDCHEM\1...\SCAN508.M (RTE Integrator)
Title : 508 scan method
Library : C:\DATABASE\WILEY7N.L

Peak Number 10 Imidazole-2,4,5-d3 \$\$ 1H-Im... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.99	0.05 ug/L	31235	Acenaphthene-d10	18.34

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Imidazole-2,4,5-d3 \$\$ 1H-Imidazo...	71	C3HD3N2	006745-43-3	5
2			Imidazole-2,4,5-d3 \$\$ 1H-Imidazo...	71	C3HD3N2	006745-43-3	4
3			Cyanic acid, ethyl ester \$\$ Ethy...	71	C3H5NO	000627-48-5	3
4			Cyanic acid, ethyl ester (CAS) \$...	71	C3H5NO	000627-48-5	3

NO
GOOD
MATCH



Library Search Compound Report

Data File : C:\MSDCHEM\1\5973N\02JAN24\00168.D
 Acq On : 24 Jan 2002 2:59 pm
 Sample : 508scan
 Misc : January 24, 2002
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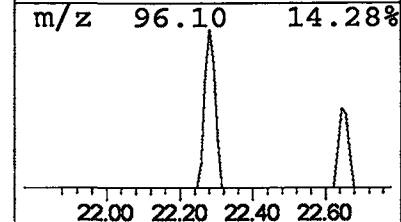
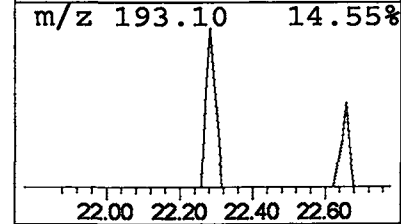
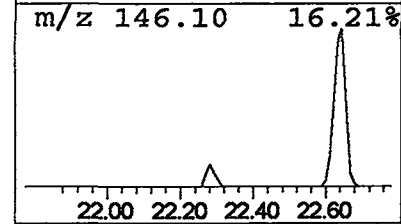
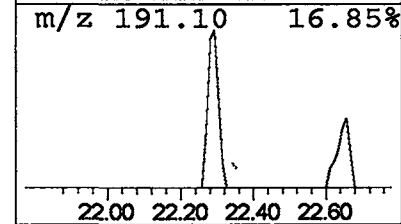
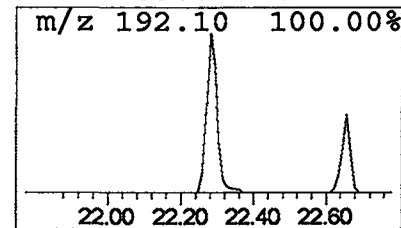
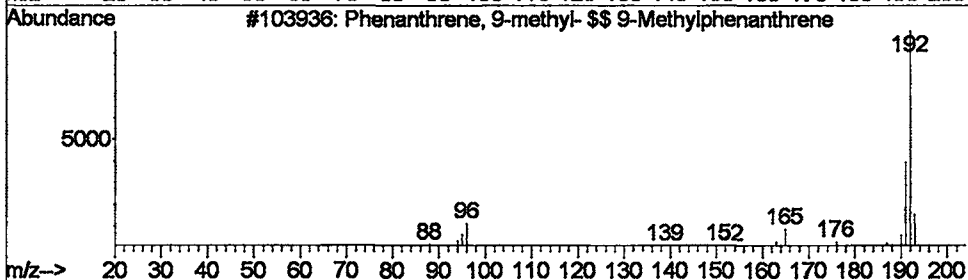
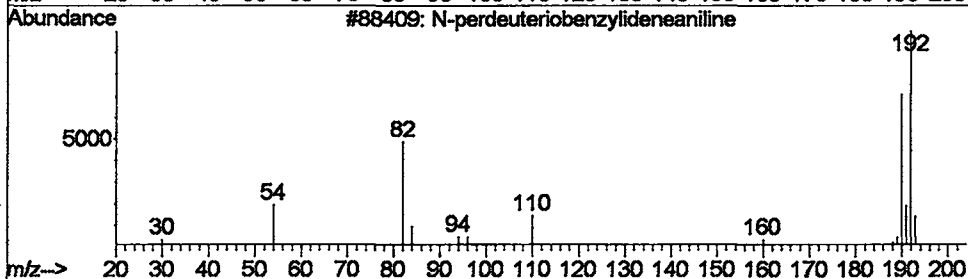
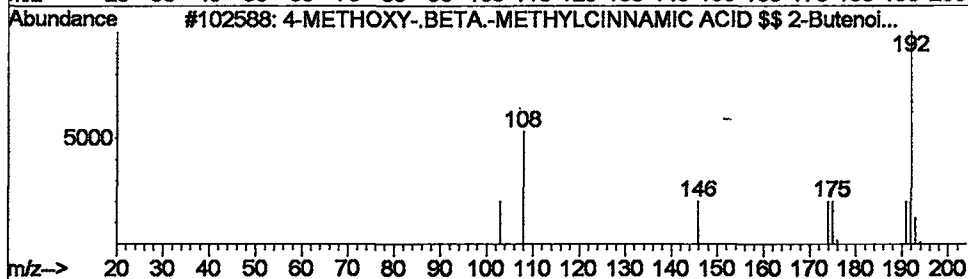
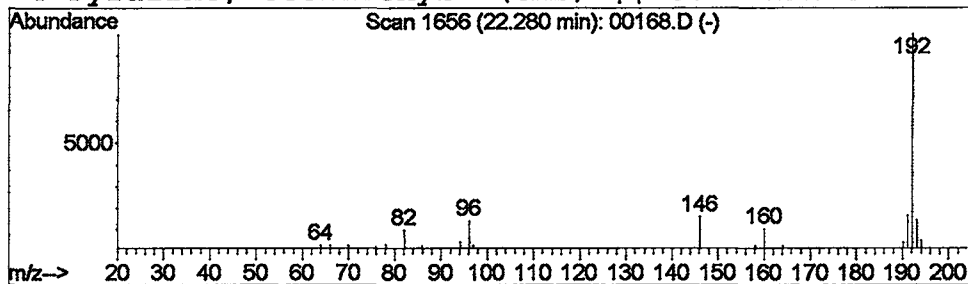
Vial: 1
 Operator:
 Inst : Instrumen
 Multiplr: 1.00

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

 Peak Number 11 4-METHOXY-.BETA.-METHYLCINN... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.28	0.14 ug/L	101559	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	58
3			Phenanthrene, 9-methyl- \$\$ 9-Met...	192	C15H12	000883-20-5	53
4			Pyrazine, tetraethyl- (CAS) \$\$ t...	192	C12H20N2	038325-19-8	50



Library Search Compound Report

Data File : C:\MSDCHEM\1\5973N\02JAN24\00168.D
Acq On : 24 Jan 2002 2:59 pm
Sample : 508scan
Misc : January 24, 2002
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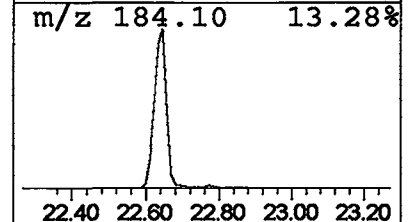
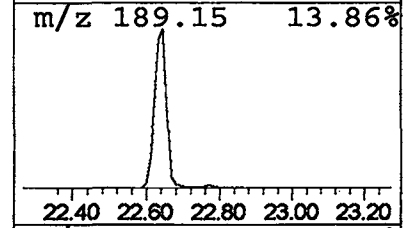
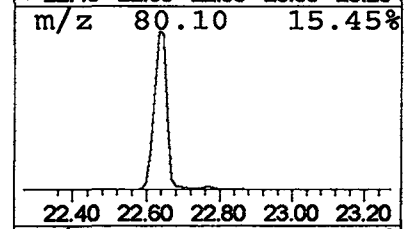
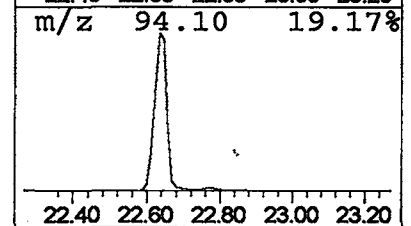
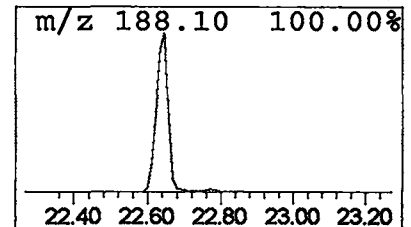
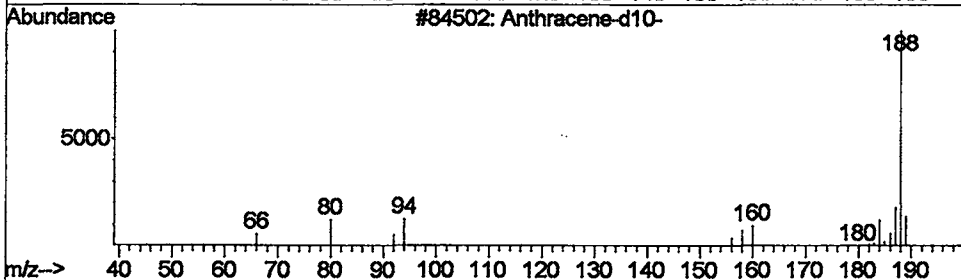
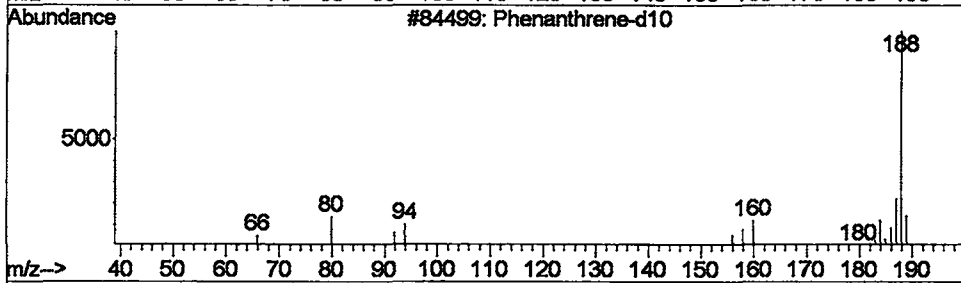
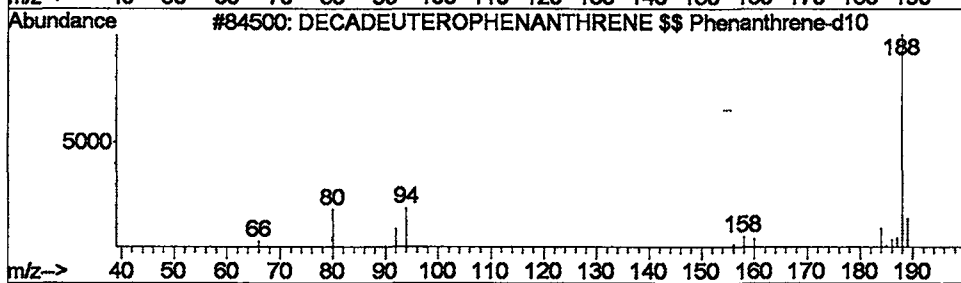
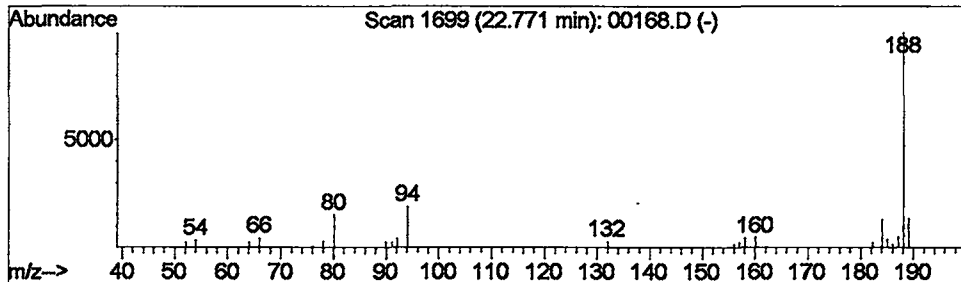
Vial: 1
Operator:
Inst : Instrumen
Multiplr: 1.00

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

Peak Number 12 DECADEUTEROPHENANTHRENE \$\$... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.77	0.32 ug/L	234038	Phenanthrene-d10	22.65

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			DECADEUTEROPHENANTHRENE \$\$ Phena...	188	C14D10	001517-22-2	91
2			Phenanthrene-d10	188	C14D10	001517-22-2	72
3			Anthracene-d10-	188	C14D10	001719-06-8	64
4			Acridine-D9	188	C13D9N	000000-00-0	59



Tentatively Identified Compound (LSC) summary

Operator ID: Date Acquired: 24 Jan 2002 2:59 pm
 Data File: C:\MSDCHEM\1\5973N\02JAN24\00168.D
 Name: 508scan
 Date: January 24, 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
Acetonitrile, dic...	3.48	0.1	ug/L	43334	1	9.72	8505730	20.0
Butanoic acid, me...	3.60	0.1	ug/L	48845	1	9.72	8505730	20.0
Phenophene, 2,3-di...	4.24	0.1	ug/L	34250	1	9.72	8505730	20.0
Methyl 1-Methyl-2...	4.62	0.1	ug/L	40152	1	9.72	8505730	20.0
3,4-Dihydropyran ...	4.85	0.2	ug/L	95892	1	9.72	8505730	20.0
Butenol, methyl-	5.01	0.1	ug/L	45260	1	9.72	8505730	20.0
Acetic acid, 3-[1...	5.89	3.8	ug/L	1597410	1	9.72	8505730	20.0
1,2-BIS (TRIMETHYL...	12.54	0.0	ug/L	23493	2	13.19	12392300	20.0
1-HYDROXY-BICYCLO...	16.55	0.0	ug/L	30979	3	18.34	13563200	20.0
Imidazole-2,4,5-d...	19.99	0.0	ug/L	31235	3	18.34	13563200	20.0
1-METHOXY-.BETA.-...	22.28	0.1	ug/L	101559	4	22.65	14490000	20.0
DECADEUTEROPHENAN...	22.77	0.3	ug/L	234038	4	22.65	14490000	20.0