

User: Vinci, David L
Westchester Dept. of Labs & Research
Date: 02/07/2002 Time: 12:52:42

Sample: AE00106
Analysis: \$GCMS GCMS Scan For Unknowns
Units: ug
Started: 2/7/02 12:52 Ended: 2/7/02 12:52 Analyst: DLV
Page: 1

	Component Name	Viol	Result	Qualifier	MDL
1	Other unknown compounds		.		
2	Surr_2,4-DDD		84		5.0

Comments for Sample AE00106 - Analysis \$GCMS

Westchester Dept. of Labs & Research

User: Vinci, David L

Date: 02-07-2002 Time: 12:53:34

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

Data File : C:\MSDCHEM\1\5973N\02JAN25\00106.D

Vial: 5

Acq On : 25 Jan 2002 1:28 pm

Operator:

Sample : 508 scan

Inst : Instrumen

Misc : January 25, 2002

Multiplr: 1.00

MS Integration Params: SPLIT.P

Quant Time: Jan 28 09:00:02 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	1557744	20.00	ug/L	0.00
10) Naphthalene-d8	13.19	136	6513174	20.00	ug/L	-0.02
17) Acenaphthene-d10	18.34	164	3321943	20.00	ug/L	-0.01
29) Phenanthrene-d10	22.65	188	5480358	20.00	ug/L	-0.02
38) Chrysene-d12	30.50	240	4873133	20.00	ug/L	-0.01
44) Perylene-d12	34.42	264	3440877	20.00	ug/L	0.00

System Monitoring Compounds

37) 2,4-DDD	27.73	235	362437	4.18	ug/L	0.00
Spiked Amount	5.000		Recovery	=	83.60%	

Target Compounds

						Qvalue
20) Dimethylphthalate	18.34	163	530504	2.42	ug/L	# 57
21) 2,6-Dinitrotoluene	18.34	165	420452	9.83	ug/L	# 17
42) Bis(2-ethylhexyl)phthalate	31.09	149	14715	0.61	ug/L	84

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

00106.D SCAN508.M Mon Jan 28 09:00:03 2002

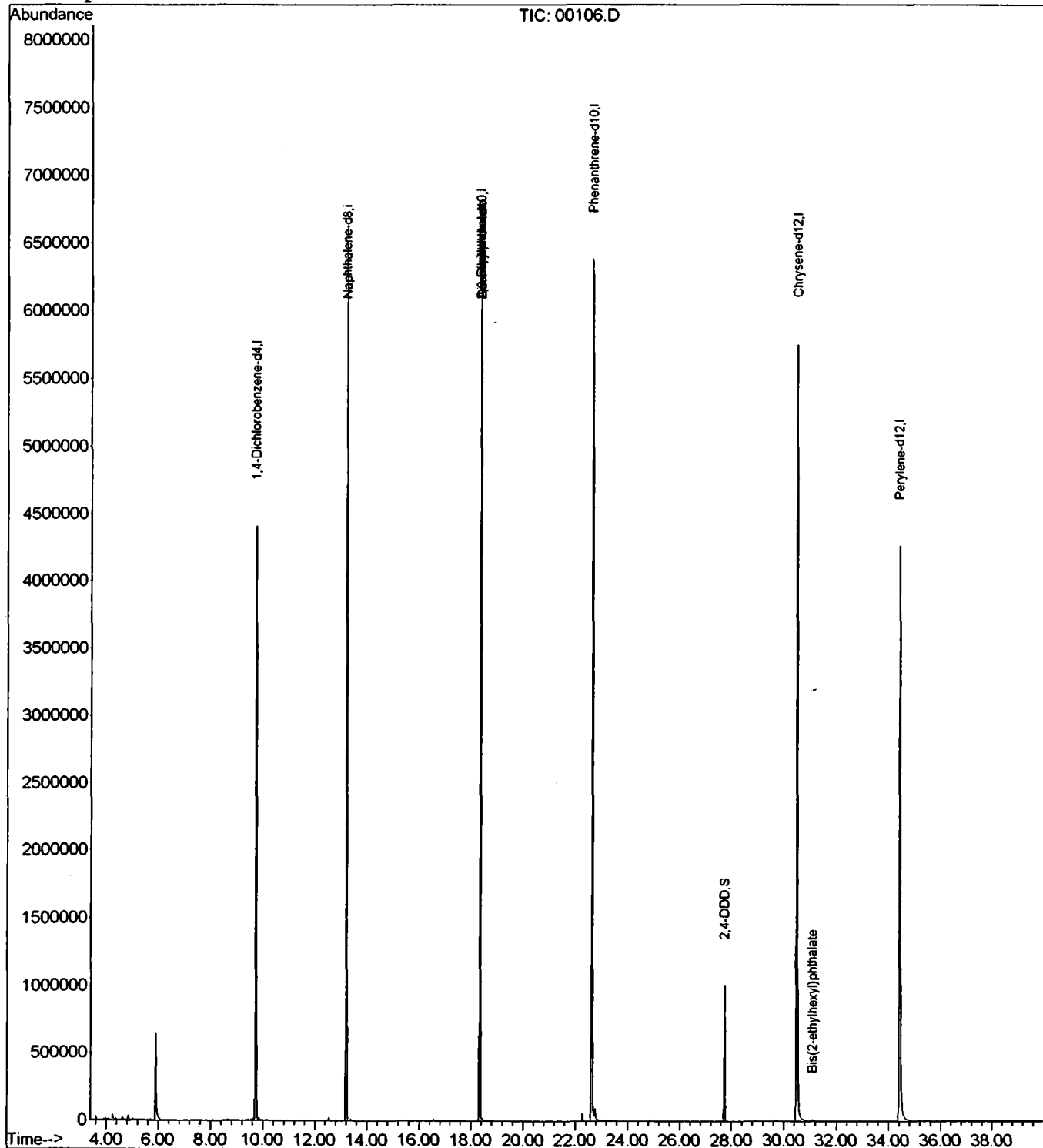
Page 1

Data File : C:\MSDCHEM\1\5973N\02JAN25\00106.D
 Acq On : 25 Jan 2002 1:28 pm
 Sample : 508 scan
 Misc : January 25, 2002
 MS Integration Params: SPLIT.P
 Quant Time: Jan 28 9:00 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



Data File : C:\MSDCHEM\1\5973N\02JAN25\00106.D
 Acq On : 25 Jan 2002 1:28 pm
 Sample : 508 scan
 Misc : January 25, 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

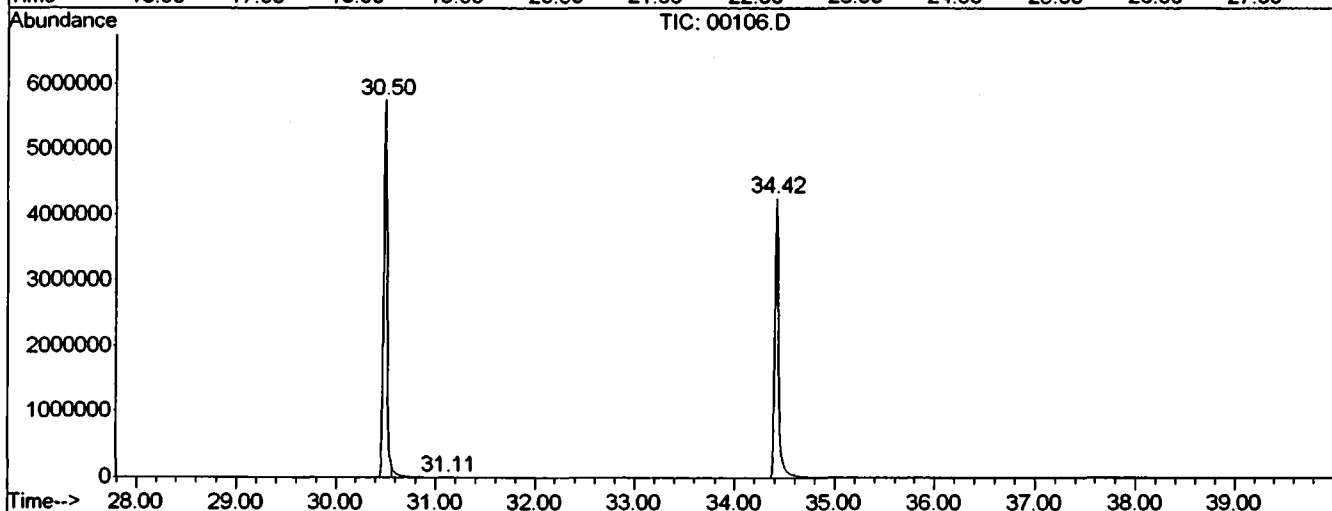
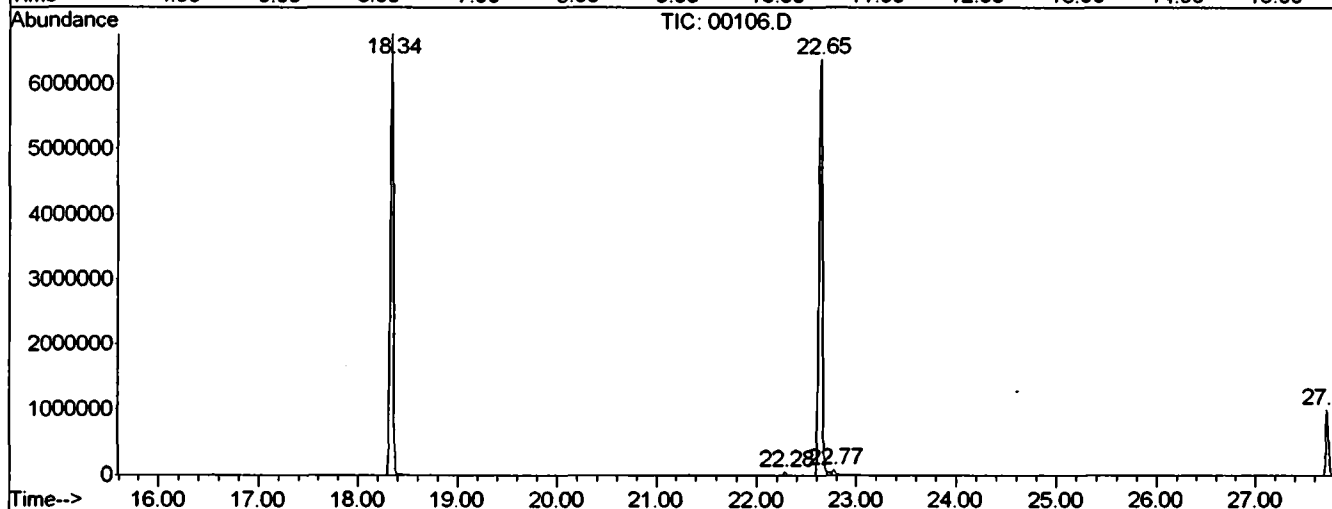
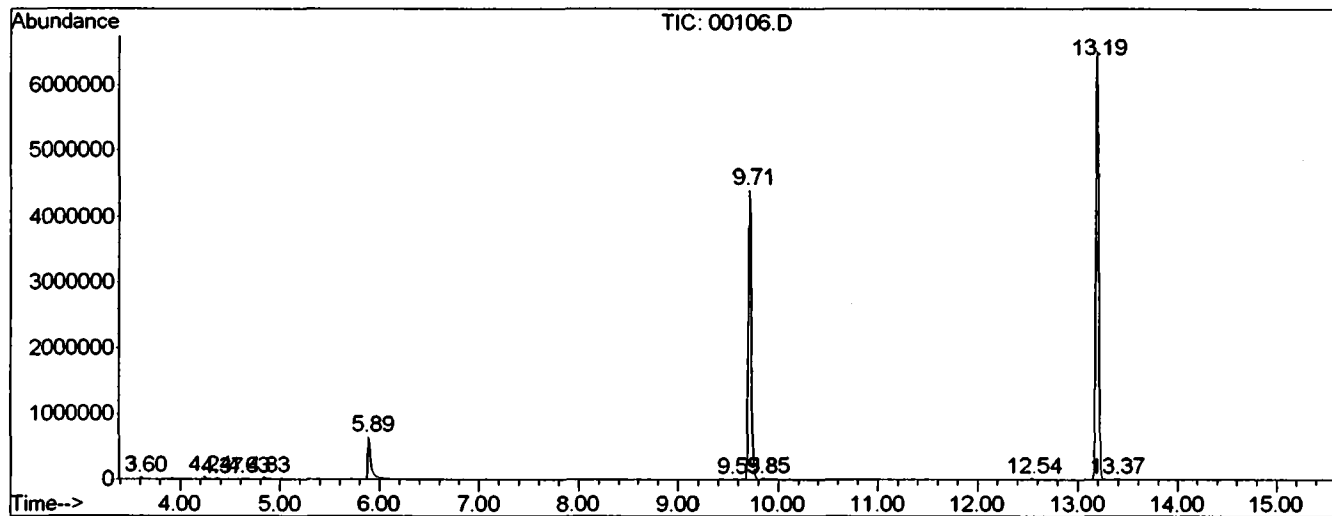
Signal : TIC

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.602	17	20	25	rBV	29725	50856	0.35%	0.064%
2	4.241	71	76	83	rVV3	42289	87635	0.60%	0.110%
3	4.367	83	87	91	rVB3	16639	36636	0.25%	0.046%
4	4.629	107	110	117	rVB3	19126	43655	0.30%	0.055%
5	4.835	123	128	141	rBV3	33516	88598	0.61%	0.112%
6	5.885	217	220	243	rBV	649470	1489126	10.21%	1.874%
7	9.550	537	541	547	rBV3	12112	32446	0.22%	0.041%
8	9.710	551	555	563	rVB	4404223	8915759	61.15%	11.222%
9	9.847	563	567	573	rBV	18743	50822	0.35%	0.064%
10	12.541	799	803	811	rVB	25876	44730	0.31%	0.056%
11	13.192	855	860	865	rBV	6511424	12579349	86.27%	15.833%
12	13.375	873	876	883	rVB2	14373	31979	0.22%	0.040%
13	18.341	1305	1311	1315	rBV	6758882	13477411	92.43%	16.963%
14	22.280	1651	1656	1661	rVB	59491	114694	0.79%	0.144%
15	22.645	1681	1688	1693	rBV2	6382840	14581203	100.00%	18.353%
16	22.771	1697	1699	1717	rVB	88460	223435	1.53%	0.281%
17	27.726	2129	2133	2139	rVV	1005721	1806287	12.39%	2.273%
18	30.500	2369	2376	2381	rBV	5753710	14007326	96.06%	17.630%
19	31.105	2419	2429	2437	rBV7	13959	57691	0.40%	0.073%
20	34.416	2713	2719	2749	rBV	4262981	11730304	80.45%	14.764%

Sum of corrected areas: 79449942

LSC Report - Integrated Chromatogram

File : C:\MSDCHEM\1\5973N\02JAN25\00106.D
 Operator :
 Acquired : 25 Jan 2002 1:28 pm using AcqMethod SCAN508
 Instrument : Instrumen
 Sample Name: 508 scan
 Misc Info : January 25, 2002
 Vial Number: 5
 Quant File :SCAN508.RES (RTE Integrator)



Data File : C:\MSDCHEM\1\5973N\02JAN25\00106.D
 Acq On : 25 Jan 2002 1:28 pm
 Sample : 508 scan
 Misc : January 25, 2002
 MS Integration Params: LSCINT.P

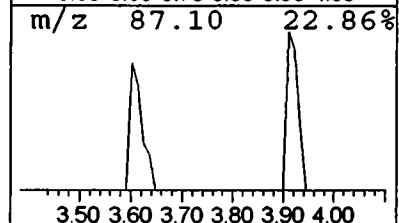
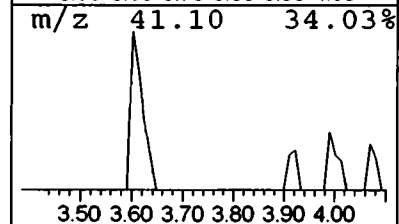
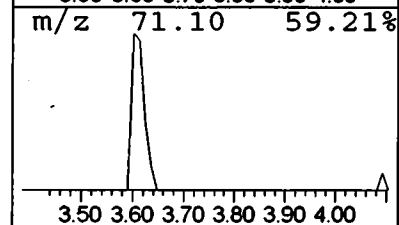
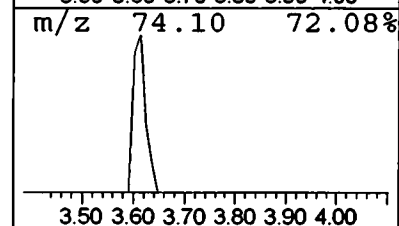
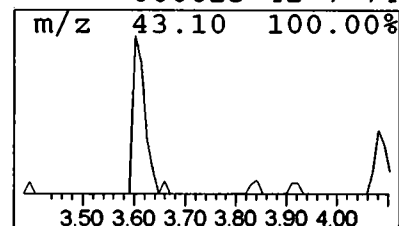
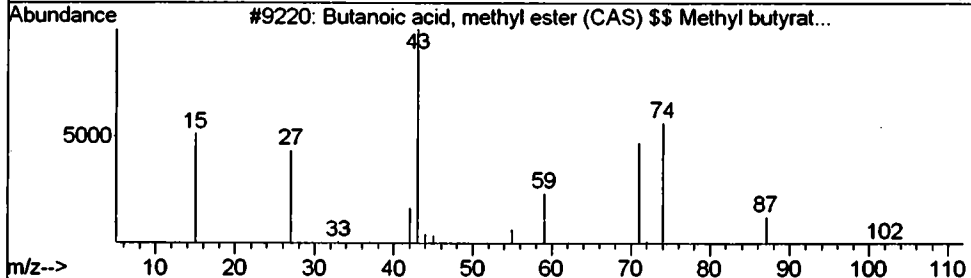
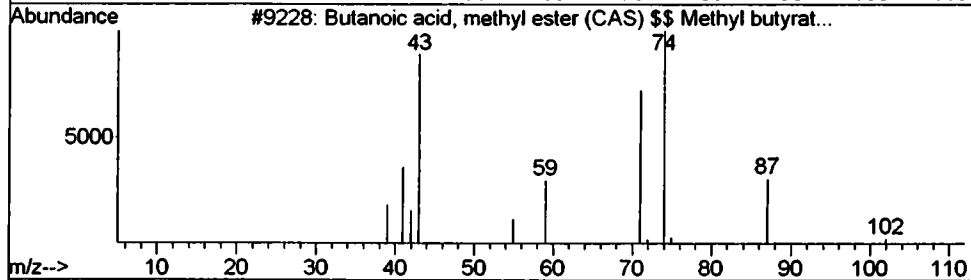
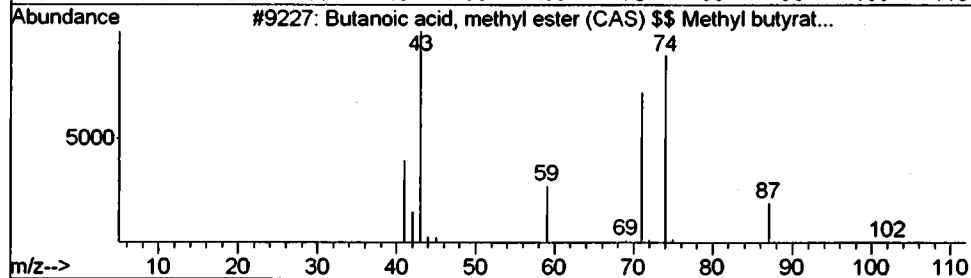
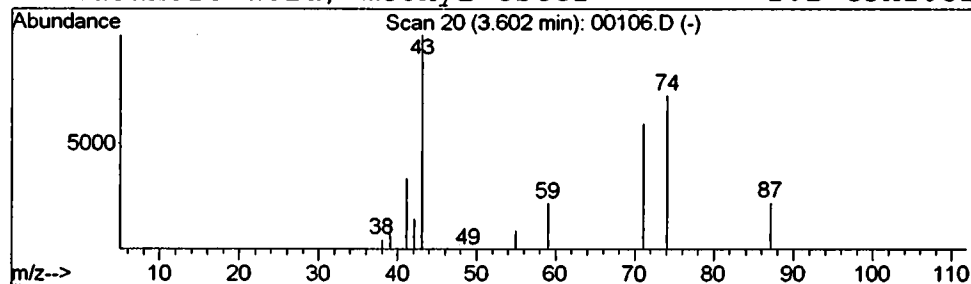
Vial: 5
 Operator:
 Inst : Instrumen
 Multiplr: 1.00

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

 Peak Number 1 Butanoic acid, methyl ester... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.60	0.11 ug/L	50856	1,4-Dichlorobenzene-d4	9.71

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	78
3		Butanoic acid, methyl ester (CAS...	102	C5H10O2	000623-42-7	78
4		Butanoic acid, methyl ester	102	C5H10O2	000623-42-7	74



Data File : C:\MSDCHEM\1\5973N\02JAN25\00106.D
 Acq On : 25 Jan 2002 1:28 pm
 Sample : 508 scan
 Misc : January 25, 2002
 MS Integration Params: LSCINT.P

Vial: 5
 Operator:
 Inst : Instrumen
 Multiplr: 1.00

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

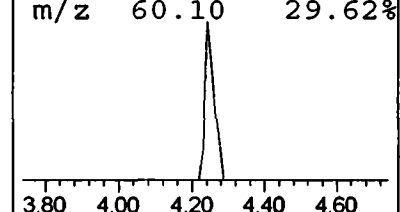
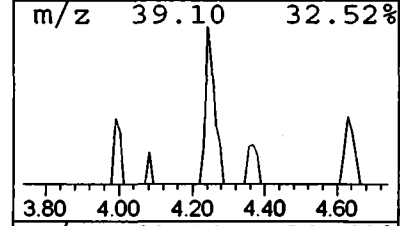
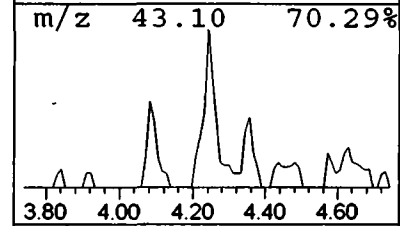
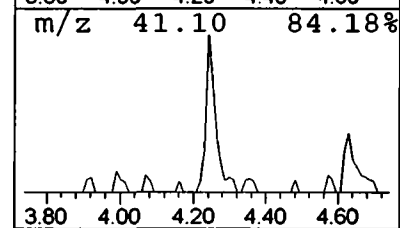
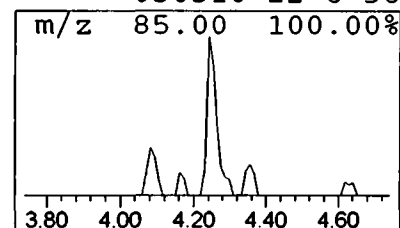
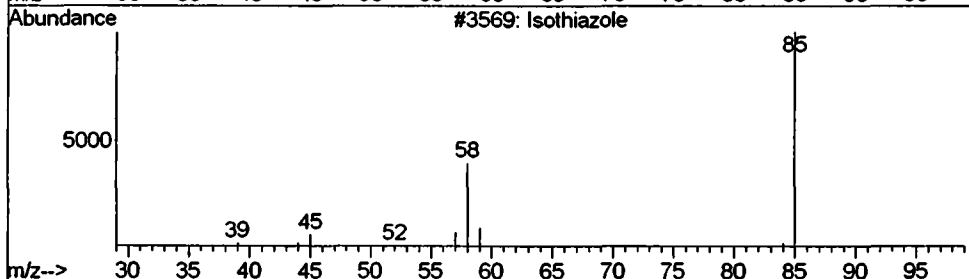
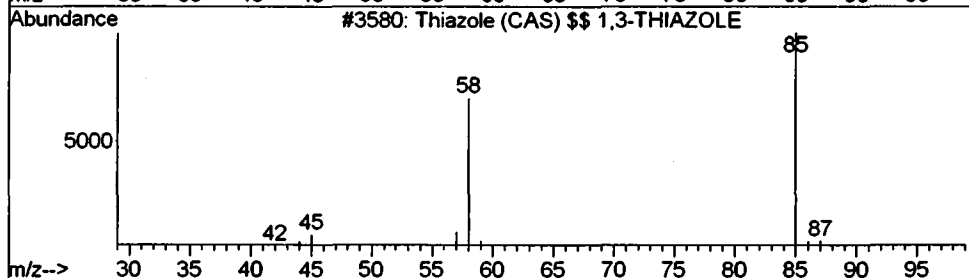
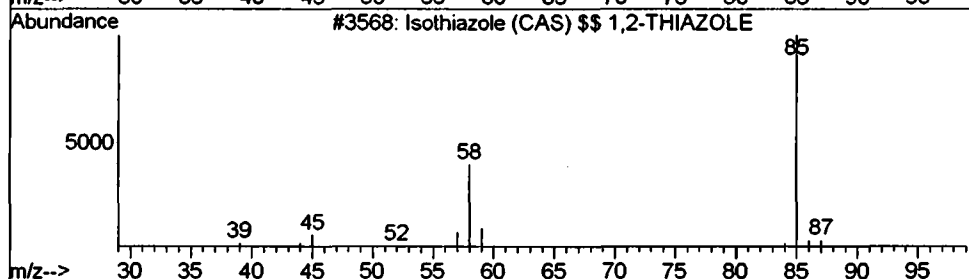
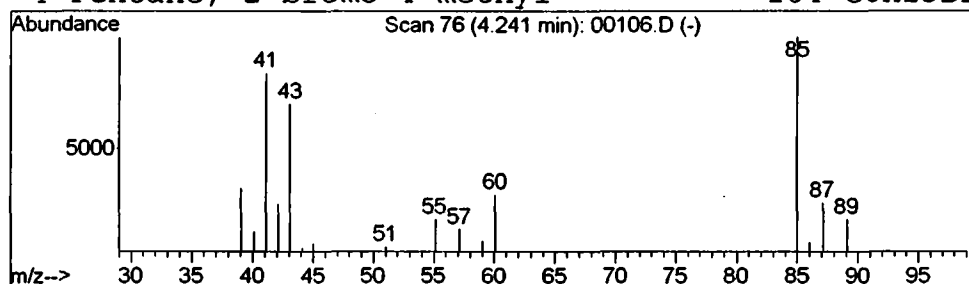
Title : 508 scan method

Library : C:\DATABASE\WILEY7N.L

 Peak Number 2 Isothiazole (CAS) \$\$ 1,2-TH... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.24	0.20 ug/L	87635	1,4-Dichlorobenzene-d4	9.71

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Isothiazole (CAS) \$\$ 1,2-THIAZOLE	85	C3H3NS	000288-16-4	43
2		Thiazole (CAS) \$\$ 1,3-THIAZOLE	85	C3H3NS	000288-47-1	38
3		Isothiazole	85	C3H3NS	000288-16-4	38
4		Pentane, 2-bromo-4-methyl-	164	C6H13Br	030310-22-6	36



Data File : C:\MSDCHEM\1\5973N\02JAN25\00106.D

Acq On : 25 Jan 2002 1:28 pm

Sample : 508 scan

Misc : January 25, 2002

MS Integration Params: LSCINT.P

Vial: 5

Operator:

Inst : Instrumen

Multiplr: 1.00

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

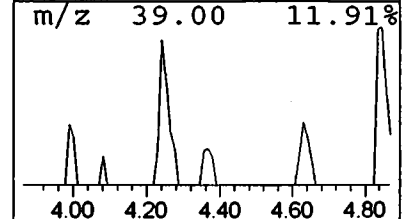
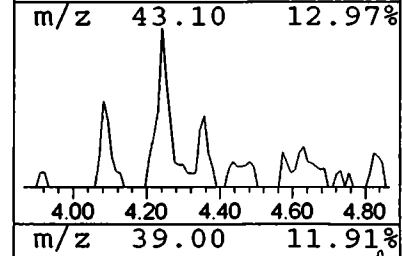
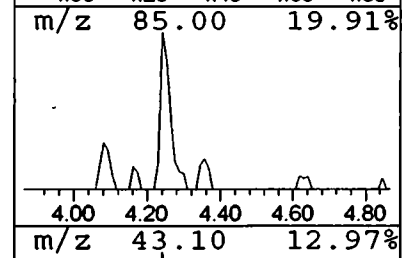
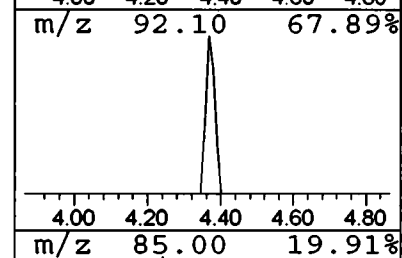
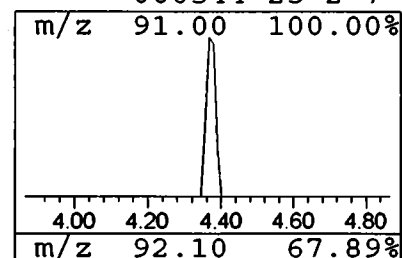
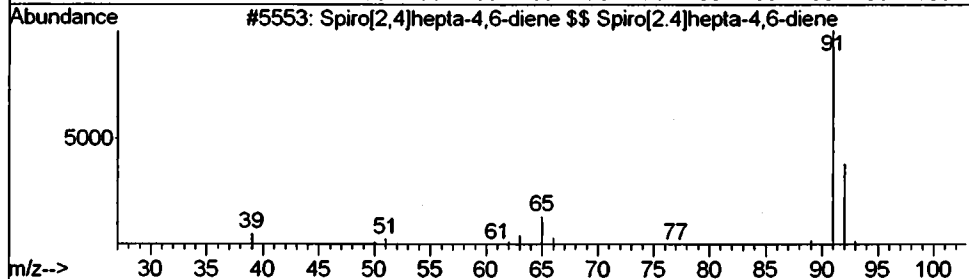
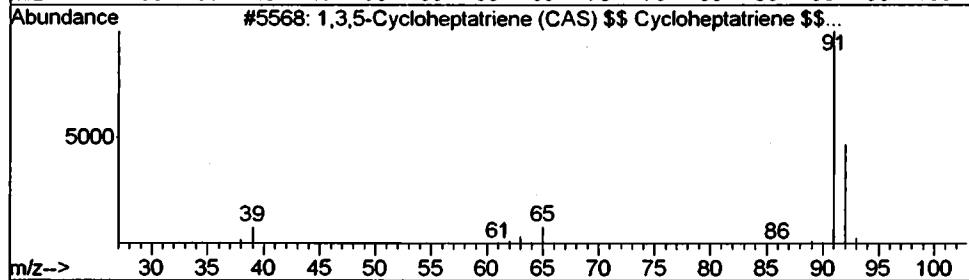
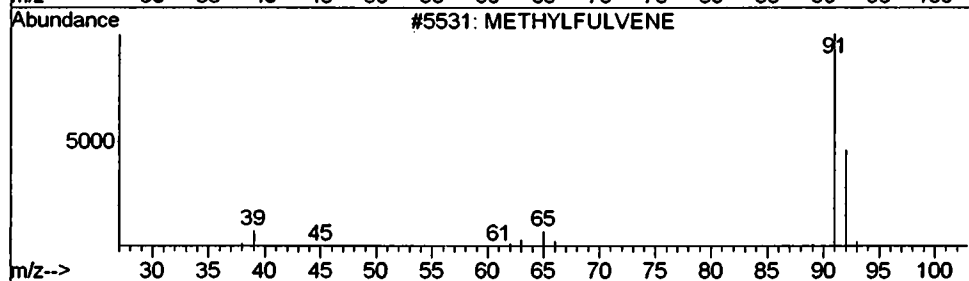
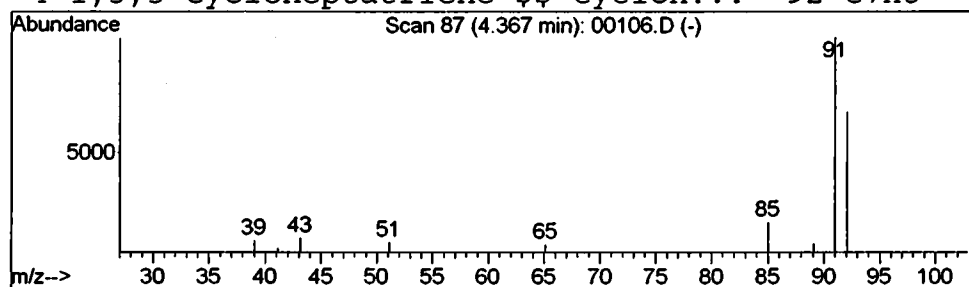
Title : 508 scan method

Library : C:\DATABASE\WILEY7N.L

Peak Number 3 METHYLFULVENE Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.37	0.08 ug/L	36636	1,4-Dichlorobenzene-d4	9.71

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		METHYLFULVENE	92	C7H8	000000-00-0	9
2		1,3,5-Cycloheptatriene (CAS) \$\$...	92	C7H8	000544-25-2	9
3		Spiro[2,4]hepta-4,6-diene \$\$ Spi...	92	C7H8	000765-46-8	7
4		1,3,5-Cycloheptatriene \$\$ Cycloh...	92	C7H8	000544-25-2	7



Data File : C:\MSDCHEM\1\5973N\02JAN25\00106.D

Acq On : 25 Jan 2002 1:28 pm

Sample : 508 scan

Misc : January 25, 2002

MS Integration Params: LSCINT.P

Vial: 5

Operator:

Inst : Instrumen

Multiplr: 1.00

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

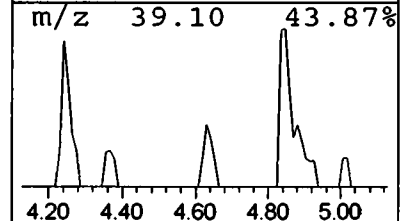
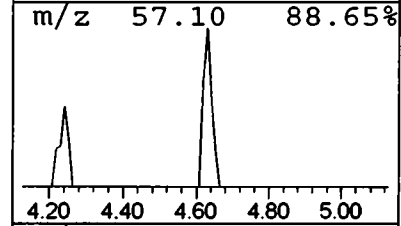
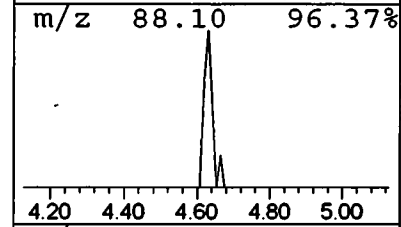
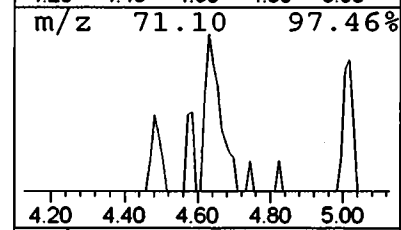
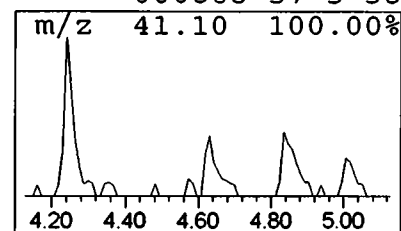
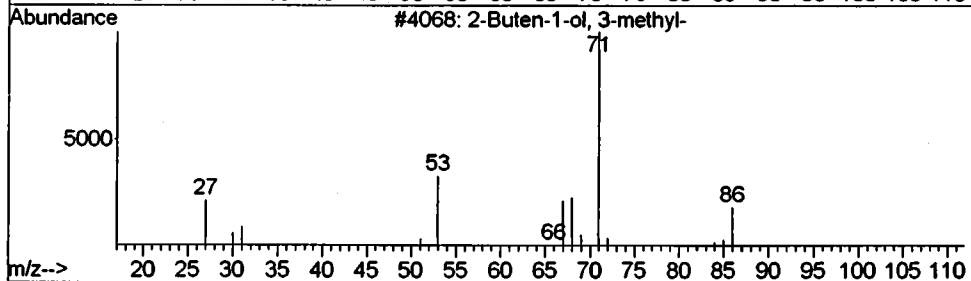
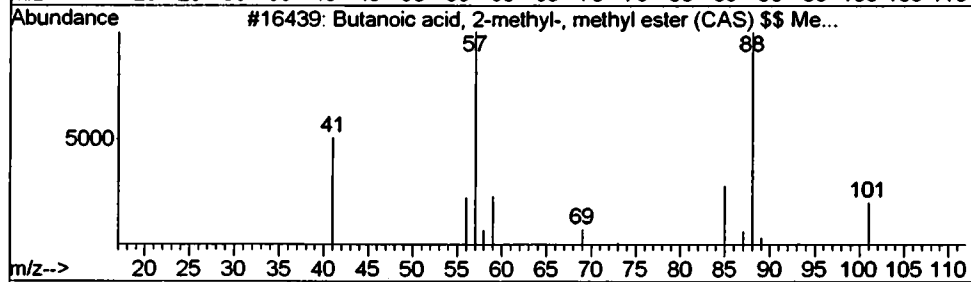
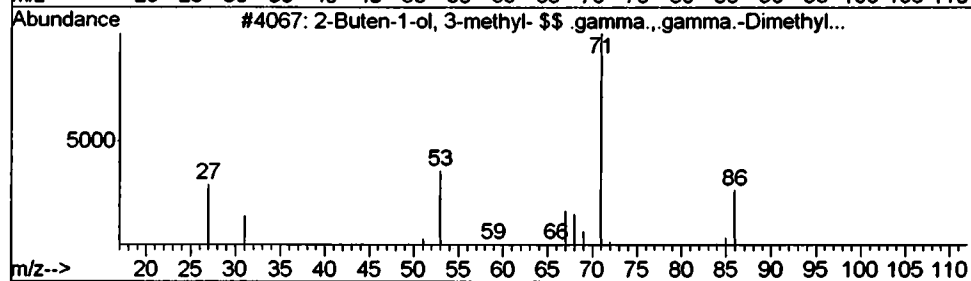
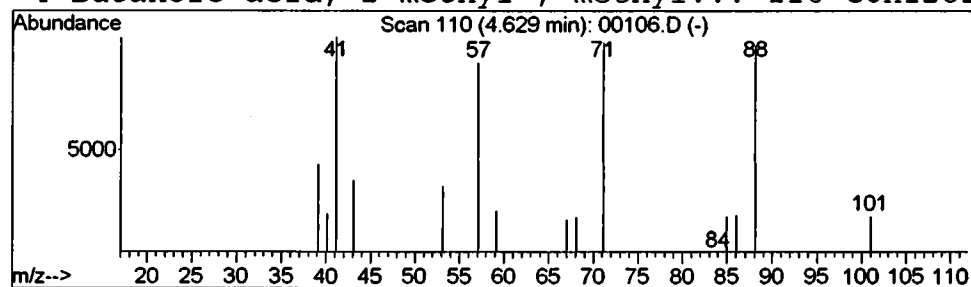
Title : 508 scan method

Library : C:\DATABASE\WILEY7N.L

Peak Number 4 2-Buten-1-ol, 3-methyl- \$\$... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.63	0.10 ug/L	43655	1,4-Dichlorobenzene-d4	9.71

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Buten-1-ol, 3-methyl- \$\$.gamm...	86	C5H10O	000556-82-1	38
2			Butanoic acid, 2-methyl-, methyl...	116	C6H12O2	000868-57-5	38
3			2-Buten-1-ol, 3-methyl-	86	C5H10O	000556-82-1	38
4			Butanoic acid, 2-methyl-, methyl...	116	C6H12O2	000868-57-5	38



Data File : C:\MSDCHEM\1\5973N\02JAN25\00106.D
 Acq On : 25 Jan 2002 1:28 pm
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 MS Integration Params: LSCINT.P

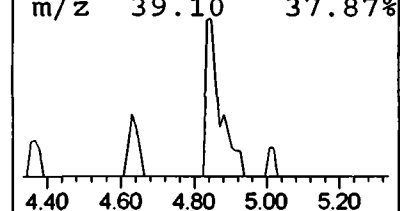
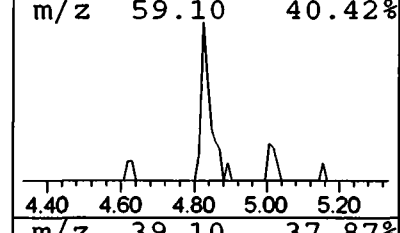
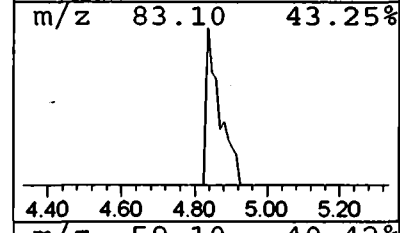
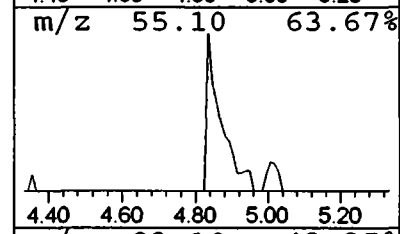
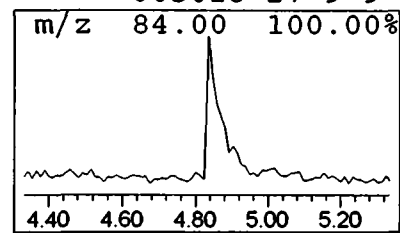
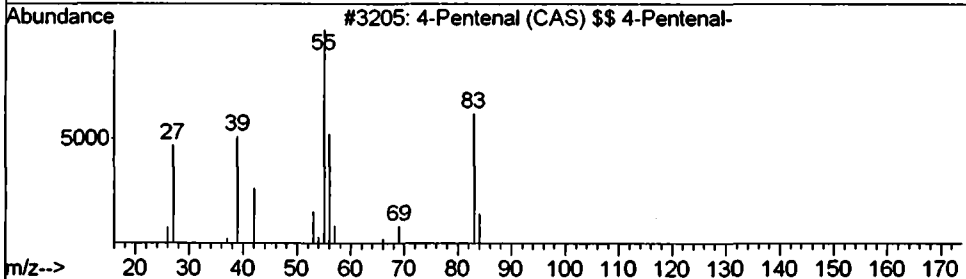
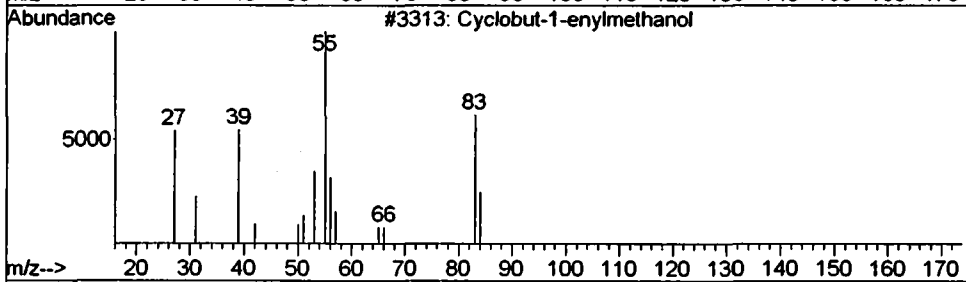
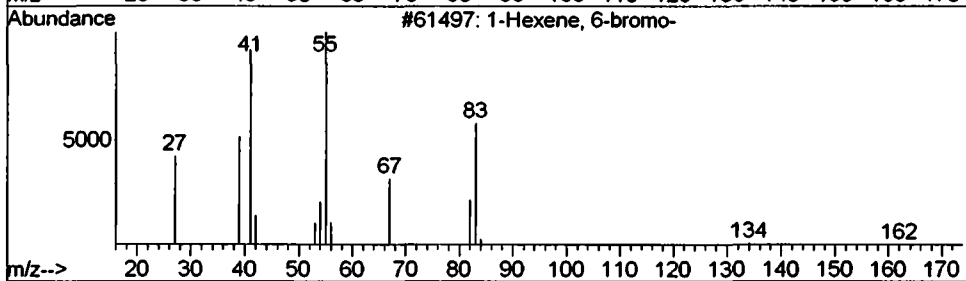
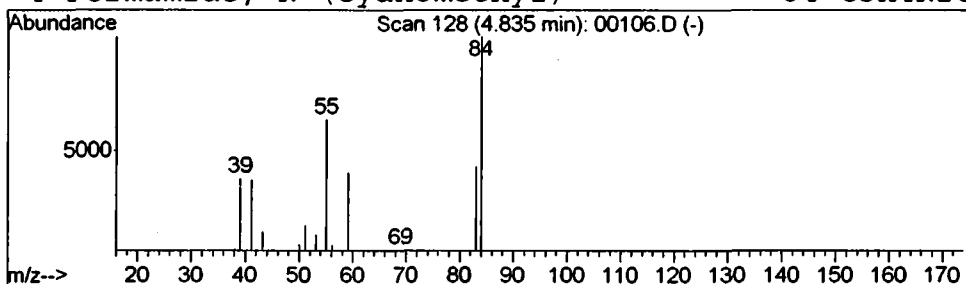
Vial: 5
 Operator:
 Inst : Instrumen
 Multiplr: 1.00

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

 Peak Number 5 1-Hexene, 6-bromo- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.83	0.20 ug/L	88598	1,4-Dichlorobenzene-d4	9.71

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1-Hexene, 6-bromo-	162	C6H11Br	002695-47-8	12
2		Cyclobut-1-enylmethanol	84	C5H8O	000000-00-0	10
3		4-Pentenal (CAS) \$\$ 4-Pentenal-	84	C5H8O	002100-17-6	10
4		Formamide, N-(cyanomethyl)-	84	C3H4N2O	005018-27-9	9



Data File : C:\MSDCHEM\1\5973N\02JAN25\00106.D
 Acq On : 25 Jan 2002 1:28 pm
 Sample : 508 scan
 Misc : January 25, 2002
 MS Integration Params: LSCINT.P

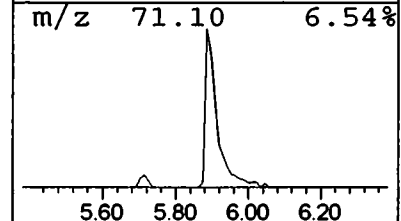
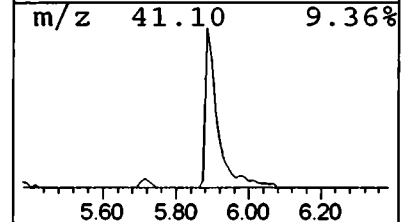
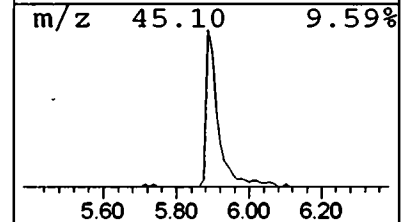
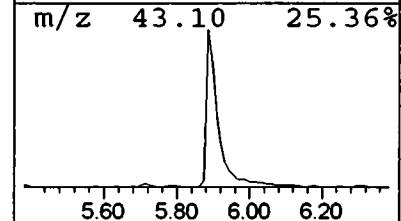
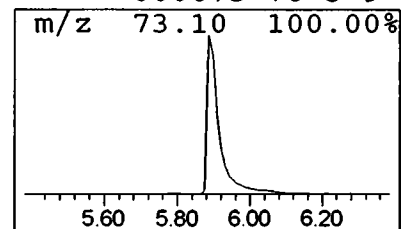
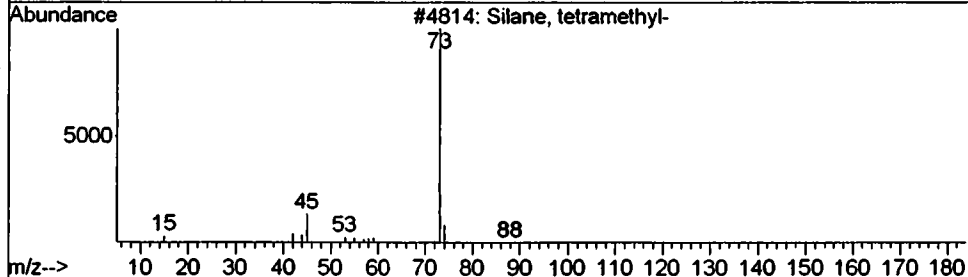
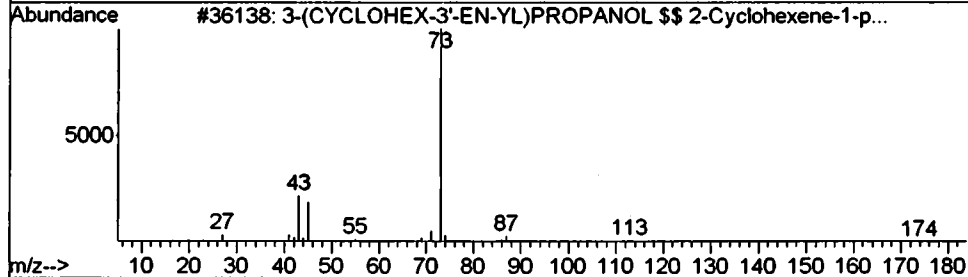
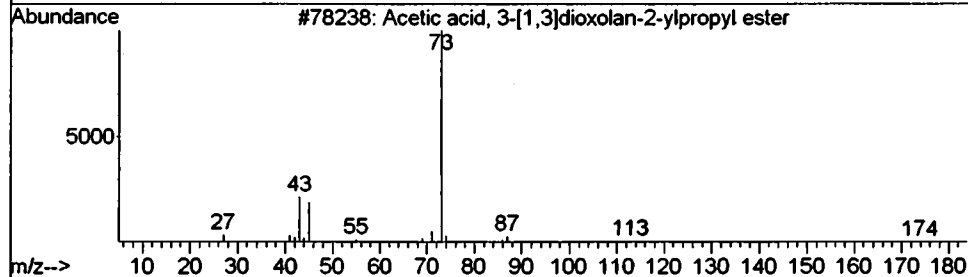
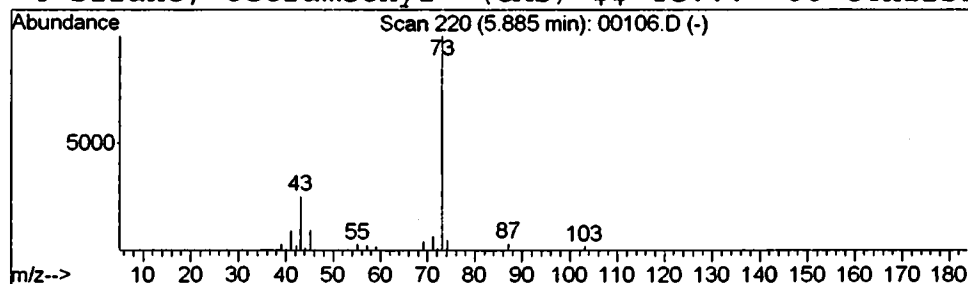
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 Operator:
 Inst : Instrumen
 Multiplr: 1.00

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.89	3.34 ug/L	1489130	1,4-Dichlorobenzene-d4	9.71

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-	88	C4H12Si	000075-76-3	32
4			Silane, tetramethyl- (CAS) \$\$ Te...	88	C4H12Si	000075-76-3	9



Data File : C:\MSDCHEM\1\5973N\02JAN25\00106.D
 Acq On : 25 Jan 2002 1:28 pm
 Sample : 508 scan
 Misc : January 25, 2002
 MS Integration Params: LSCINT.P

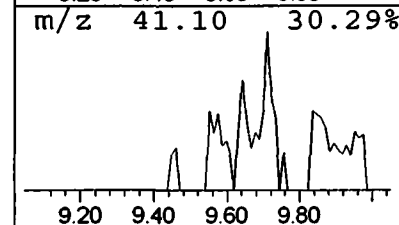
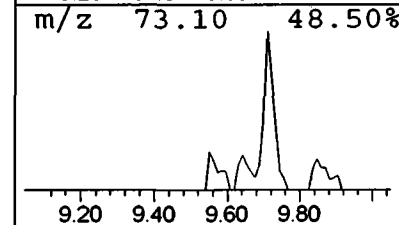
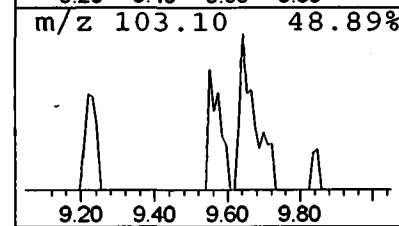
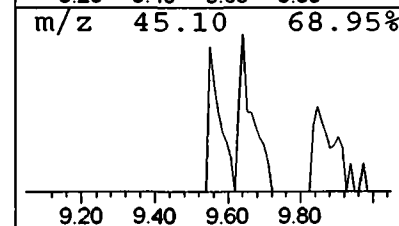
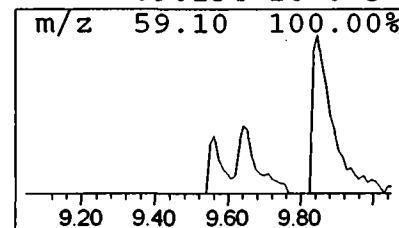
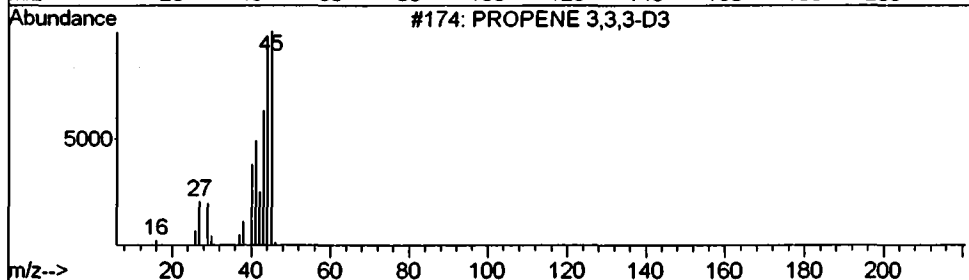
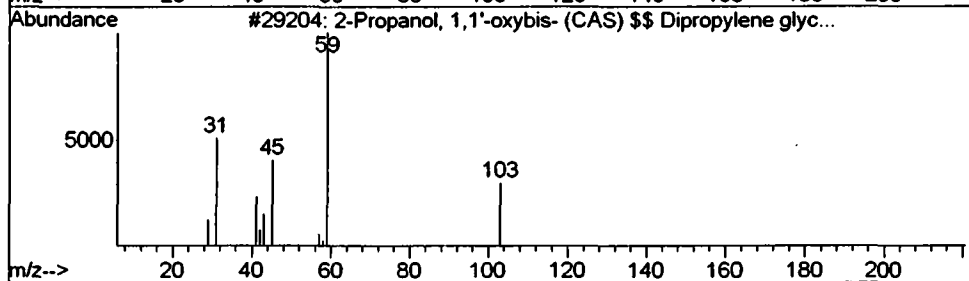
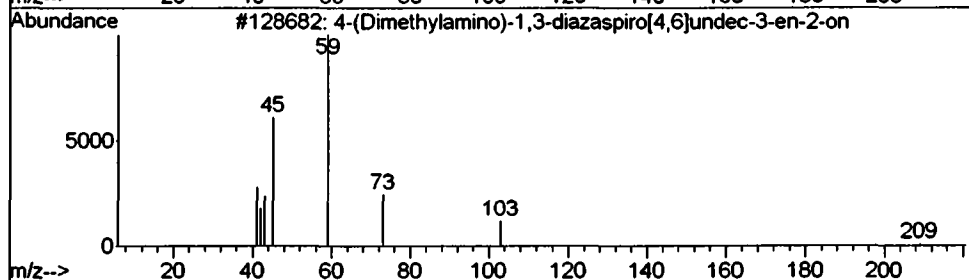
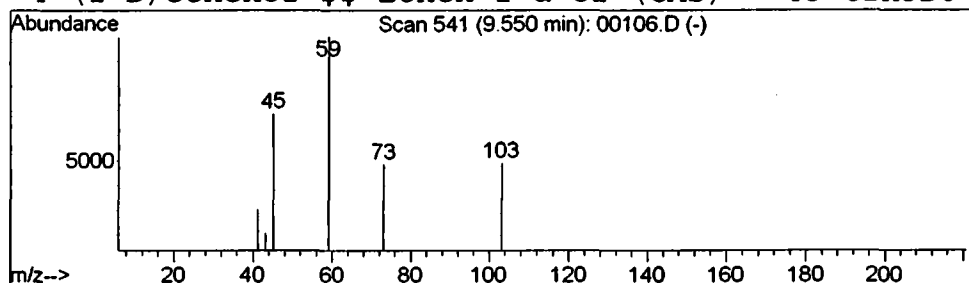
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 Inst : Instrumen
 Multiplr: 1.00

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

 Peak Number 7 4-(Dimethylamino)-1,3-diaza... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.55	0.07 ug/L	32446	1,4-Dichlorobenzene-d4	9.71

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			4-(Dimethylamino)-1,3-diazaspiro...	209	C11H19N3O	000000-00-0	74
2			2-Propanol, 1,1'-oxybis- (CAS) \$...	134	C6H14O3	000110-98-5	9
3			PROPENE 3,3,3-D3	45	C3H3D3	001517-51-7	7
4			(1-D)ethenol \$\$ Ethen-1-d-ol (CAS)	45	C2H3DO	090134-16-0	5



Data File : C:\MSDCHEM\1\5973N\02JAN25\00106.D
 Acq On : 25 Jan 2002 1:28 pm
 Sample : 508 scan
 Misc : January 25, 2002
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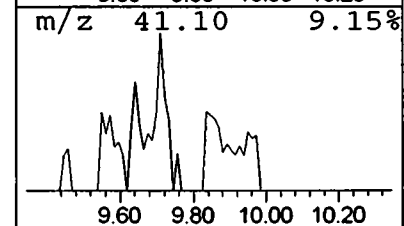
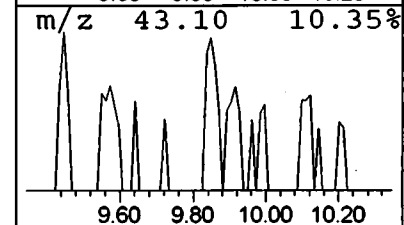
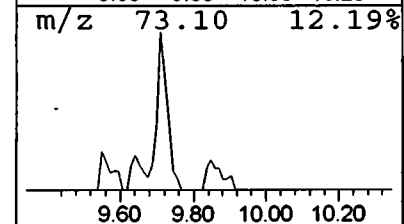
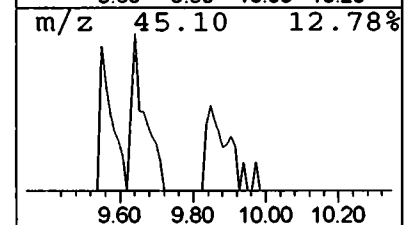
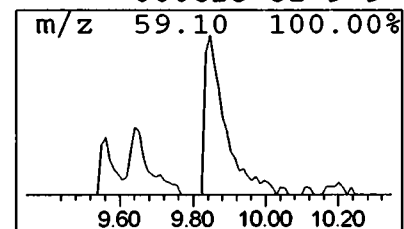
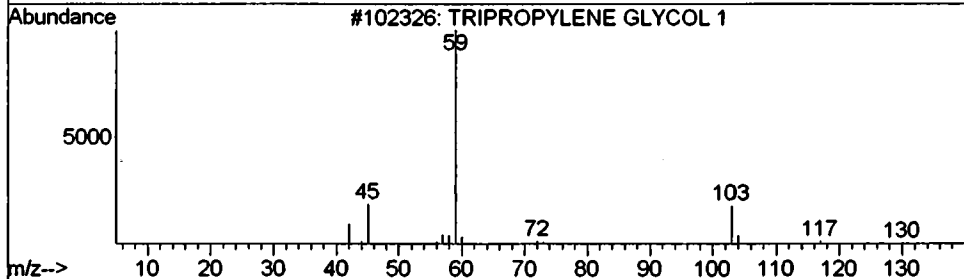
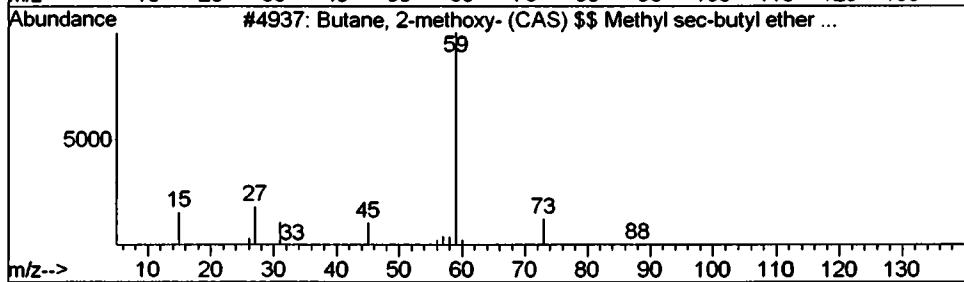
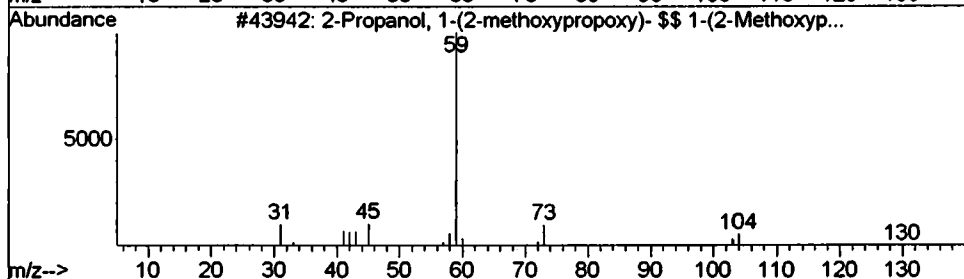
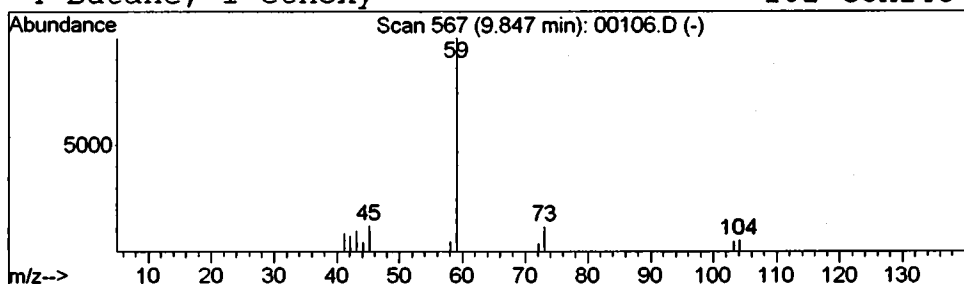
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 Inst : Instrumen
 Multiplr: 1.00

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 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.85	0.11 ug/L	50822	1,4-Dichlorobenzene-d4	9.71

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Propanol, 1-(2-methoxypropoxy)...	148	C7H16O3	013429-07-7	83
2		Butane, 2-methoxy- (CAS) \$\$ Meth...	88	C5H12O	006795-87-5	9
3		TRIPROPYLENE GLYCOL 1	192	C9H20O4	000000-00-0	9
4		Butane, 1-ethoxy-	102	C6H14O	000628-81-9	9



Data File : C:\MSDCHEM\1\5973N\02JAN25\00106.D
Acq On : 25 Jan 2002 1:28 pm
Sample : 508 scan
Misc : January 25, 2002
MS Integration Params: LSCINT.P

Vial: 5
Operator:
Inst : Instrumen
Multiplr: 1.00

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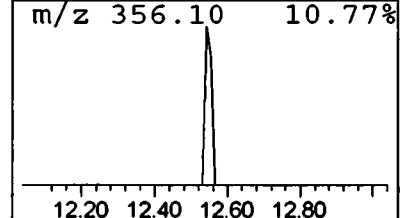
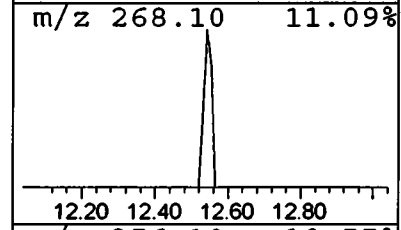
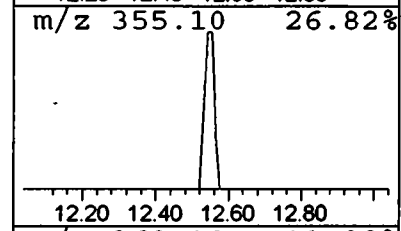
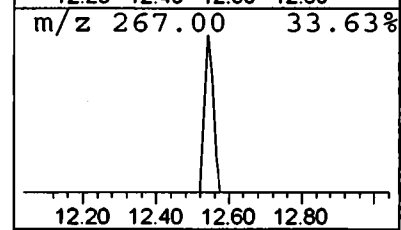
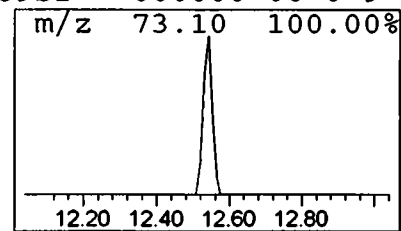
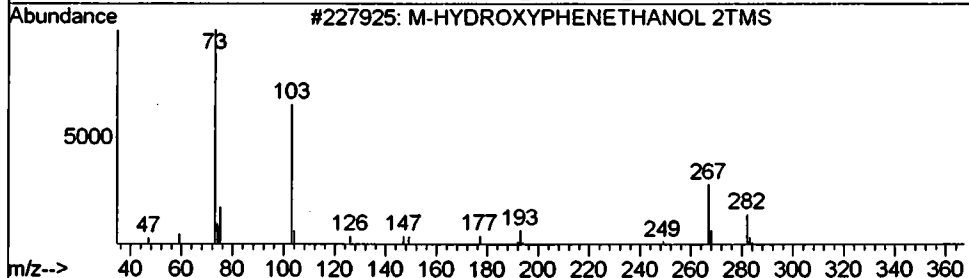
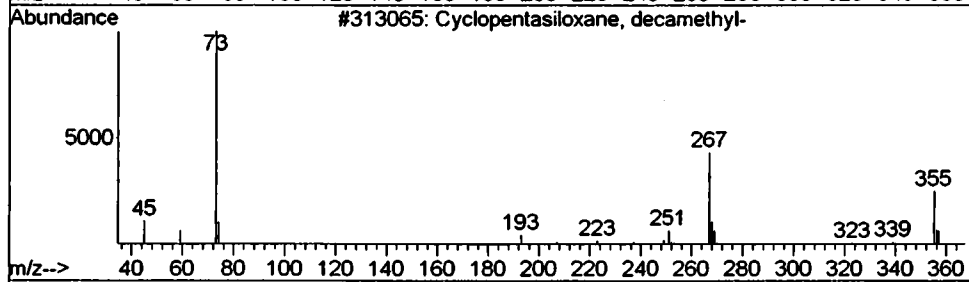
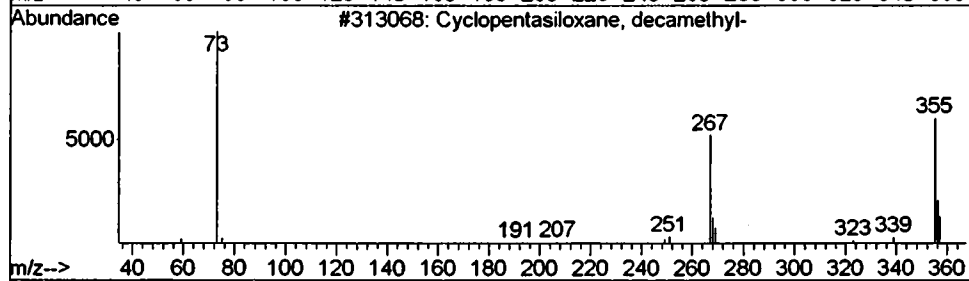
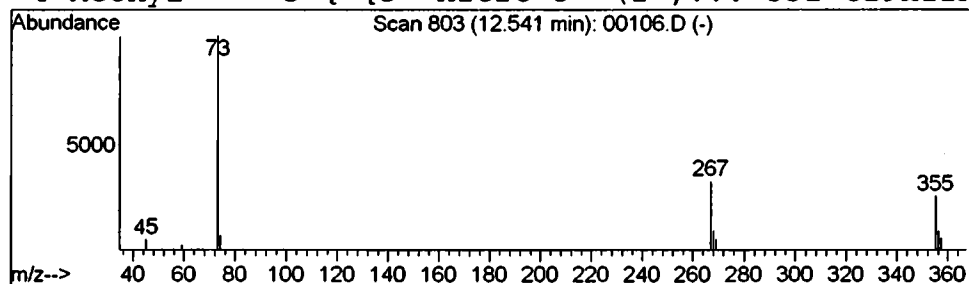
Title : 508 scan method

Library : C:\DATABASE\WILEY7N.L

Peak Number 9 Cyclopentasiloxane, decamet... Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.54	0.07 ug/L	44730	Naphthalene-d8	13.19

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclopentasiloxane, decamethyl-	370	C10H30O5Si5	000541-02-6	72
2			Cyclopentasiloxane, decamethyl-	370	C10H30O5Si5	000541-02-6	53
3			M-HYDROXYPHENETHANOL 2TMS	282	C14H26O2Si2	000000-00-0	25
4			Methyl 5-[[5'-nitro-3'-(1",...]	532	C19H12N6O9S2	000000-00-0	9



Data File : C:\MSDCHEM\1\5973N\02JAN25\00106.D
 Acq On : 25 Jan 2002 1:28 pm
 Sample : 508 scan
 Misc : January 25, 2002
 MS Integration Params: LSCINT.P

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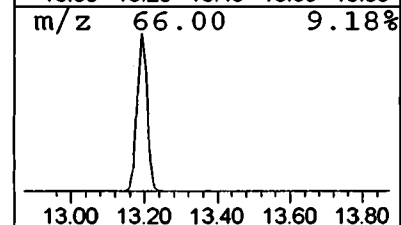
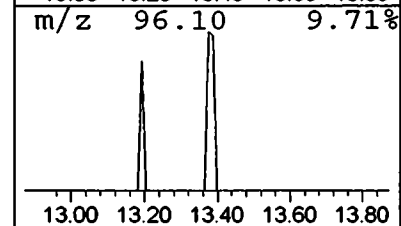
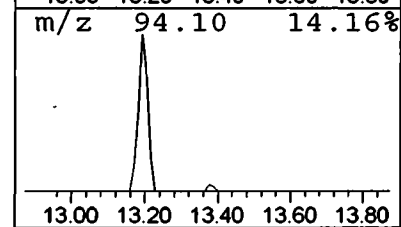
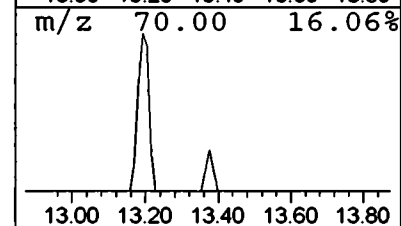
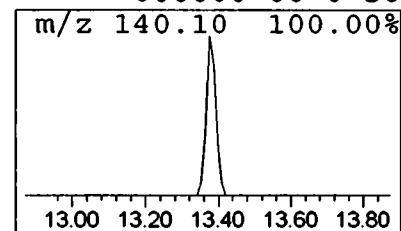
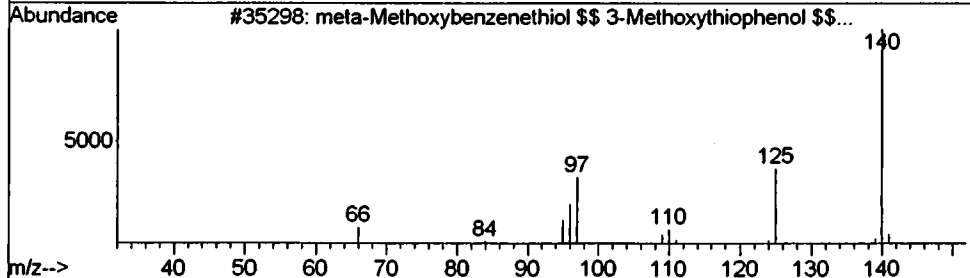
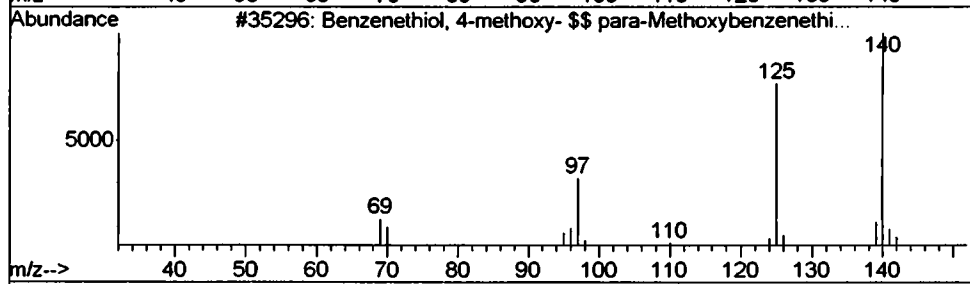
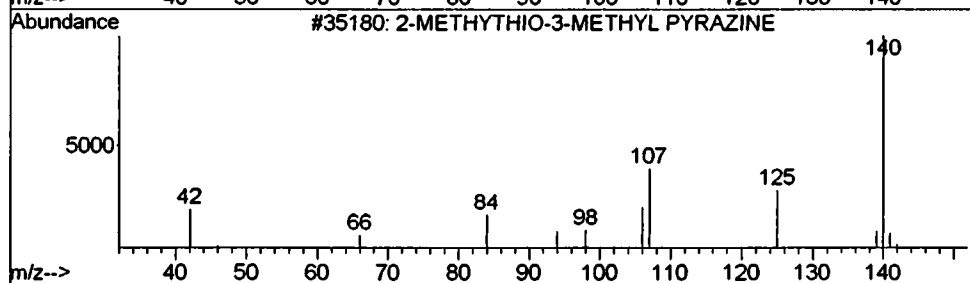
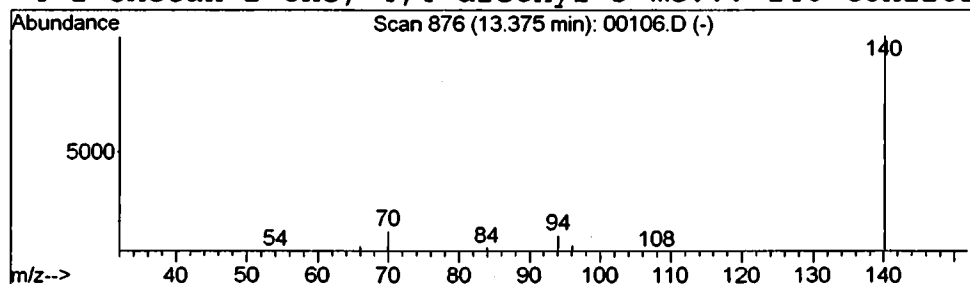
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 Title : 508 scan method
 Library : C:\DATABASE\WILEY7N.L

 Peak Number 10 2-METHYTHIO-3-METHYL PYRAZINE Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.37	0.05 ug/L	31979	Naphthalene-d8	13.19

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-METHYTHIO-3-METHYL PYRAZINE	140	C6H8N2S	002882-20-4	78
2			Benzenethiol, 4-methoxy- \$\$ para...	140	C7H8OS	000696-63-9	64
3			meta-Methoxybenzenethiol \$\$ 3-Me...	140	C7H8OS	015570-12-4	64
4			1-Oxetan-2-one, 4,4-diethyl-3-me...	140	C8H12O2	000000-00-0	56

NO
 GDM
 MATCH



Data File : C:\MSDCHEM\1\5973N\02JAN25\00106.D
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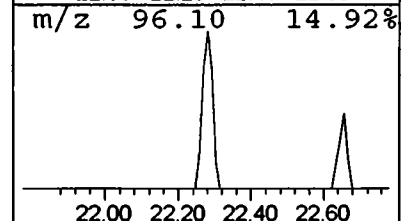
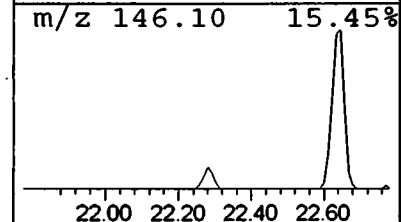
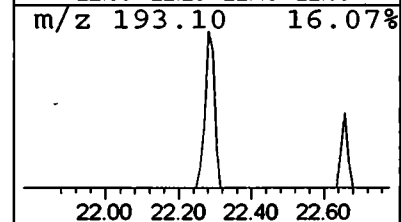
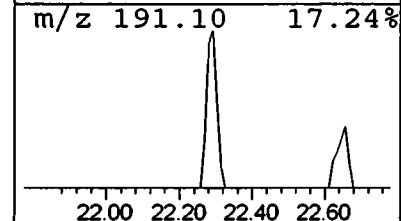
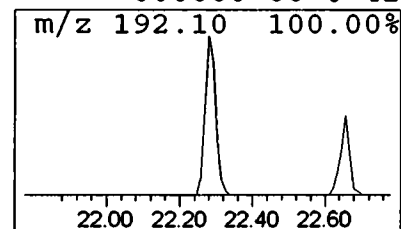
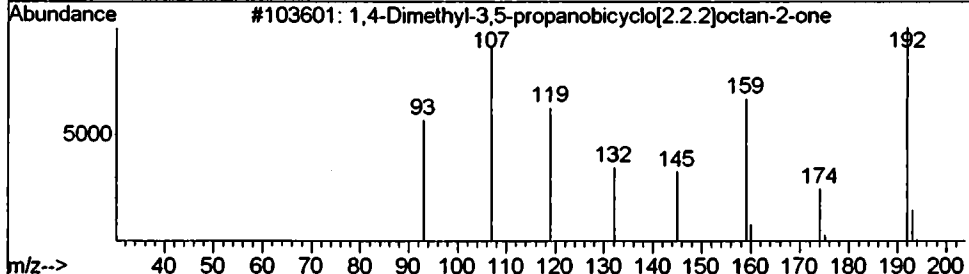
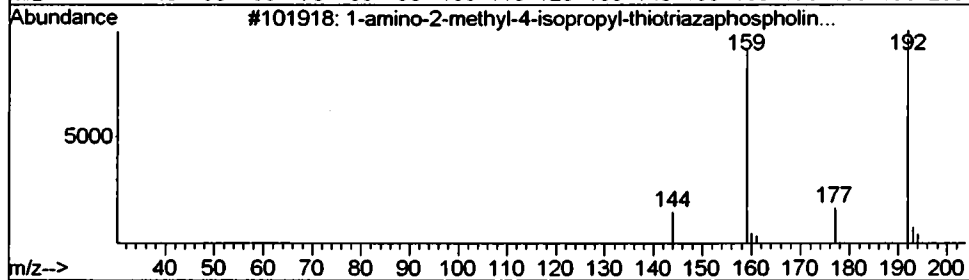
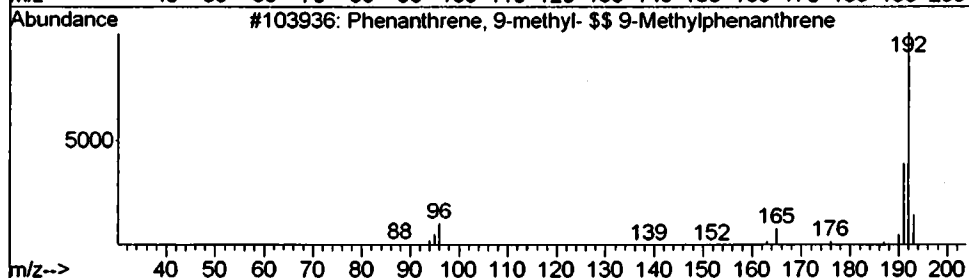
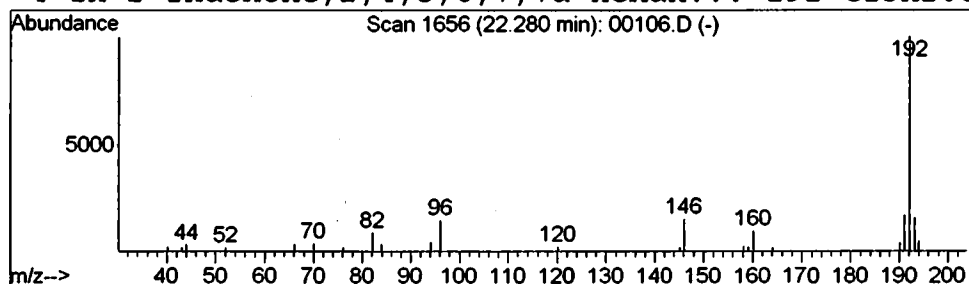
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 Operator:
 Inst : Instrumen
 Multiplr: 1.00

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

 Peak Number 11 Phenanthrene, 9-methyl- \$\$... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.28	0.16 ug/L	114694	Phenanthrene-d10	22.65

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Phenanthrene, 9-methyl- \$\$ 9-Met...	192	C15H12	000883-20-5	53
2			1-amino-2-methyl-4-isopropyl-thi...	192	C5H13N4PS	107264-54-0	50
3			1,4-Dimethyl-3,5-propanobicyclo[...	192	C13H20O	000000-00-0	47
4			1H-2-Indenone, 2,4,5,6,7,7a-hexah...	192	C13H20O	000000-00-0	42



Data File : C:\MSDCHEM\1\5973N\02JAN25\00106.D
Acq On : 25 Jan 2002 1:28 pm
Sample : 508 scan
Misc : January 25, 2002
MS Integration Params: LSCINT.P

Vial: 5
Operator:
Inst : Instrumen
Multiplr: 1.00

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

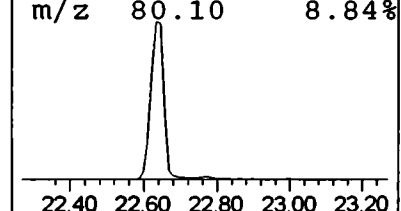
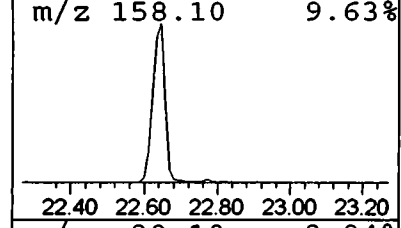
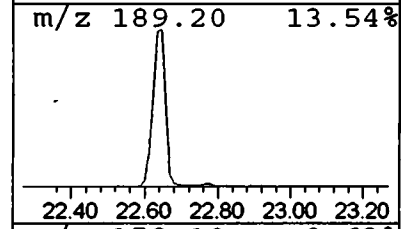
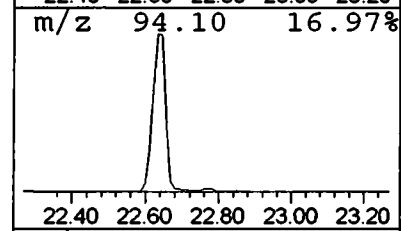
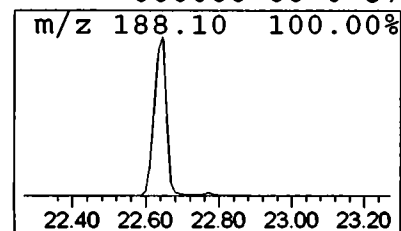
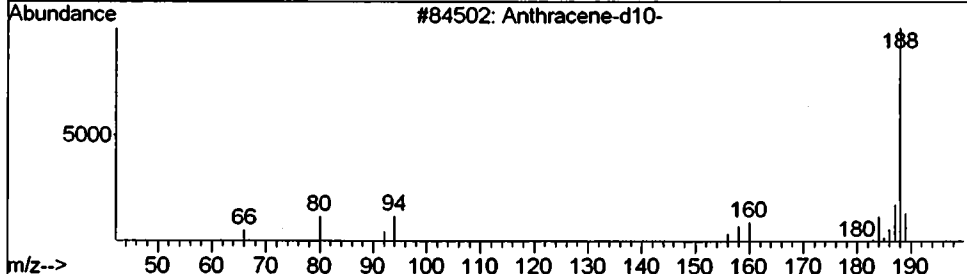
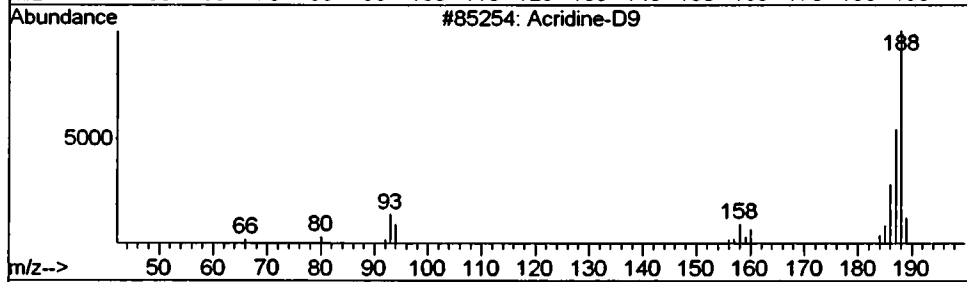
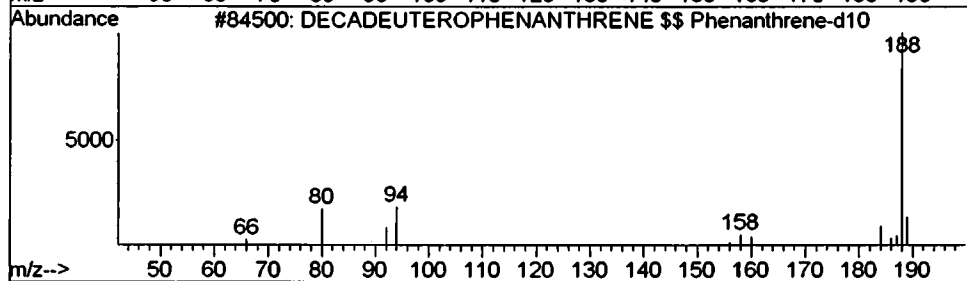
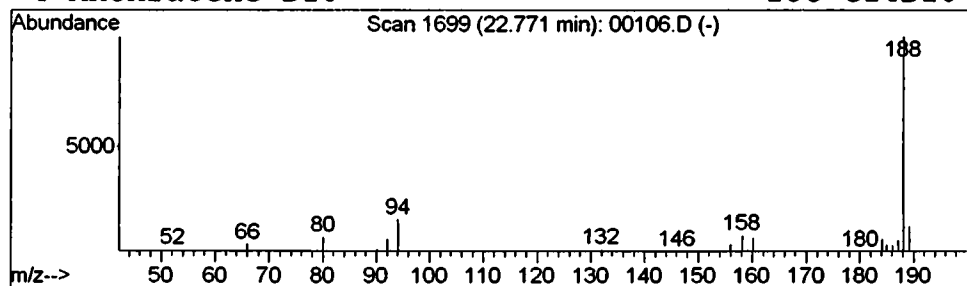
Title : 508 scan method

Library : C:\DATABASE\WILEY7N.L

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.77	0.31 ug/L	223435	Phenanthrene-d10	22.65

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			DECADEUTEROPHENANTHRENE \$\$ Phena...	188	C14D10	001517-22-2	91
2			Acridine-D9	188	C13D9N	000000-00-0	87
3			Anthracene-d10-	188	C14D10	001719-06-8	87
4			Anthracene-D10	188	C14D10	000000-00-0	87



Operator ID: Date Acquired: 25 Jan 2002 1:28 pm
Data File: C:\MSDCHEM\1\5973N\02JAN25\00106.D
Name: 508 scan
Misc: January 25, 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
Butanoic acid, me...	3.60	0.1	ug/L	50856	1	9.71	8915760	20.0
Isothiazole (CAS)...	4.24	0.2	ug/L	87635	1	9.71	8915760	20.0
METHYLFULVENE	4.37	0.1	ug/L	36636	1	9.71	8915760	20.0
2-Buten-1-ol, 3-m...	4.63	0.1	ug/L	43655	1	9.71	8915760	20.0
1-Hexene, 6-bromo-	4.83	0.2	ug/L	88598	1	9.71	8915760	20.0
Acetic acid, 3-[1...	5.89	3.3	ug/L	1489130	1	9.71	8915760	20.0
4-(Dimethylamino)...	9.55	0.1	ug/L	32446	1	9.71	8915760	20.0
2-Propanol, 1-(2-...	9.85	0.1	ug/L	50822	1	9.71	8915760	20.0
Cyclopentasiloxan...	12.54	0.1	ug/L	44730	2	13.19	12579300	20.0
2-METHYTHIO-3-MET...	13.37	0.1	ug/L	31979	2	13.19	12579300	20.0
Phenanthrene, 9-m...	22.28	0.2	ug/L	114694	4	22.65	14581200	20.0
DECADEUTEROPHENAN...	22.77	0.3	ug/L	223435	4	22.65	14581200	20.0