

User: Vinci, David L
Westchester Dept. of Labs & Research
Date: 02/08/2002 Time: 14:33:50

Sample: AD29522
Analysis: \$GCMS GCMS Scan For Unknowns
Units: ug
Started: 2/8/02 14:33 Ended: 2/8/02 14:33 Analyst: DLV
Page: 1

	Component Name	Viol	Result	Qualifier	MDL
1	Other unknown compounds		.		
2	Surr_2,4-DDD		81		5.0

Comments for Sample AD29522 - Analysis \$GCMS

Westchester Dept. of Labs & Research

User: Vinci, David L

Date: 02-08-2002 Time: 14:36:56

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

The recoveries for 9 of the 43 compounds in the LCS Standard were below their acceptance range.

Quantitation Report

(QT Reviewed)

Data File : C:\MSDCHEM\1\5973N\02FEB01\29522.D

Vial: 4

Acq On : 1 Feb 2002 1:04 pm

Operator:

Sample : 508 scan

Inst : Instrumen

Misc : February 1, 2002

Multiplr: 1.00

MS Integration Params: SPLIT.P

Quant Time: Feb 04 08:34:14 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	1488882	20.00	ug/L	0.00
10) Naphthalene-d8	13.19	136	6259691	20.00	ug/L	-0.02
17) Acenaphthene-d10	18.34	164	3172062	20.00	ug/L	-0.01
29) Phenanthrene-d10	22.63	188	5155211	20.00	ug/L	-0.03
38) Chrysene-d12	30.50	240	4414019	20.00	ug/L	-0.01
44) Perylene-d12	34.42	264	3140467	20.00	ug/L	0.00

System Monitoring Compounds

37) 2,4-DDD	27.73	235	331630	4.07	ug/L	0.00
Spiked Amount	5.000		Recovery	=	81.40%	

Target Compounds

42) Bis(2-ethylhexyl)phthalate	31.09	149	15513	0.62	ug/L	Qvalue 83
--------------------------------	-------	-----	-------	------	------	--------------

*Re extraordi
sample
nice & clean!*

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

29522.D SCAN508.M

Mon Feb 04 09:34:19 2002

Page 1

Quantitation Report

(QT Reviewed)

Data File : C:\MSDCHEM\1\5973N\02FEB01\29522.D

Vial: 4

Acq On : 1 Feb 2002 1:04 pm

Operator:

Sample : 508 scan

Inst : Instrumen

Misc : February 1, 2002

Multiplr: 1.00

MS Integration Params: SPLIT.P

Quant Time: Feb 4 9:33 2002

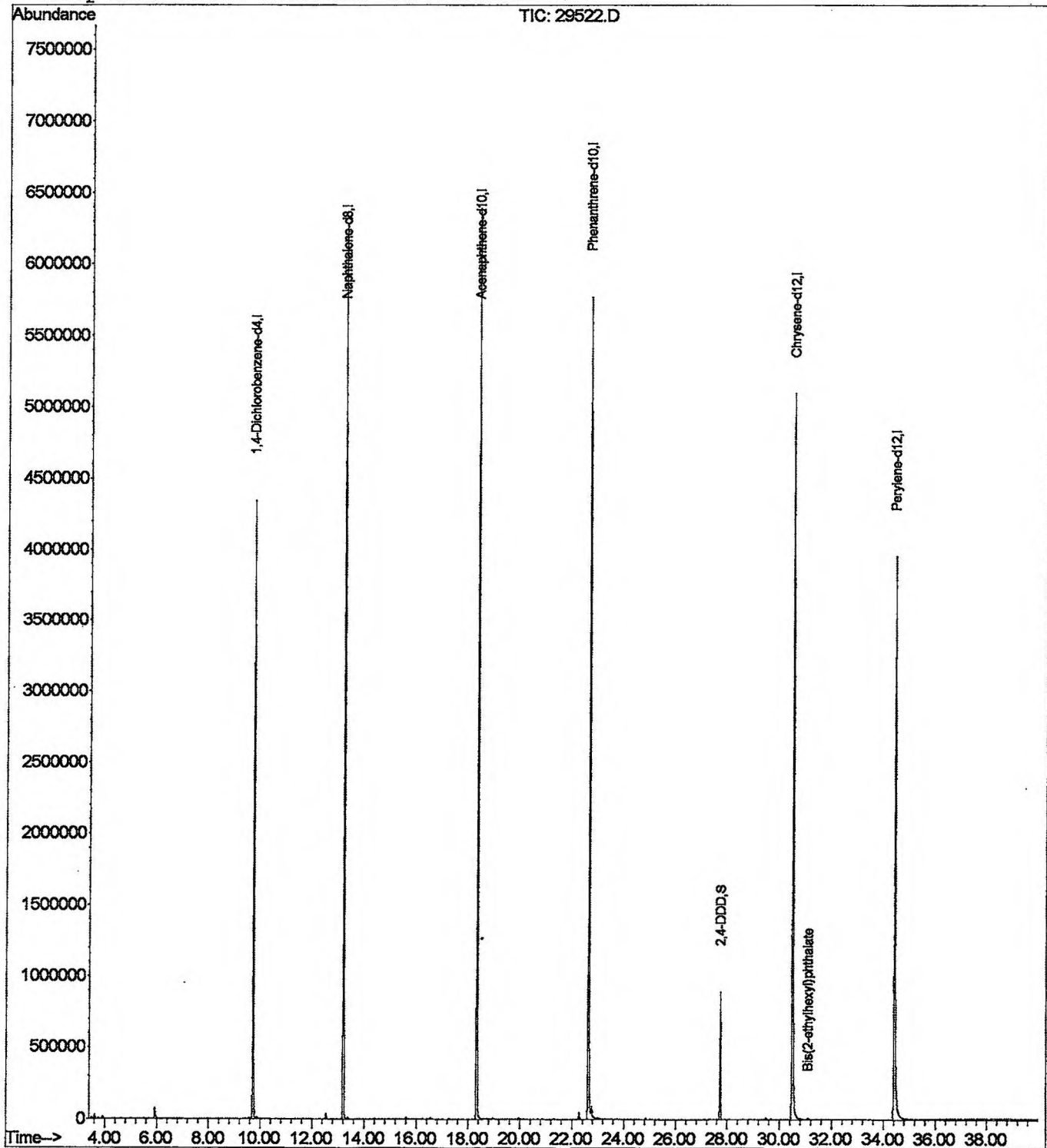
Quant Results File: SCAN508.RE

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\02FEB01\29522.D

Acq On : 1 Feb 2002 1:04 pm

Sample : 508 scan

Misc : February 1, 2002

MS Integration Params: LSCINT.P

Vial: 4

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

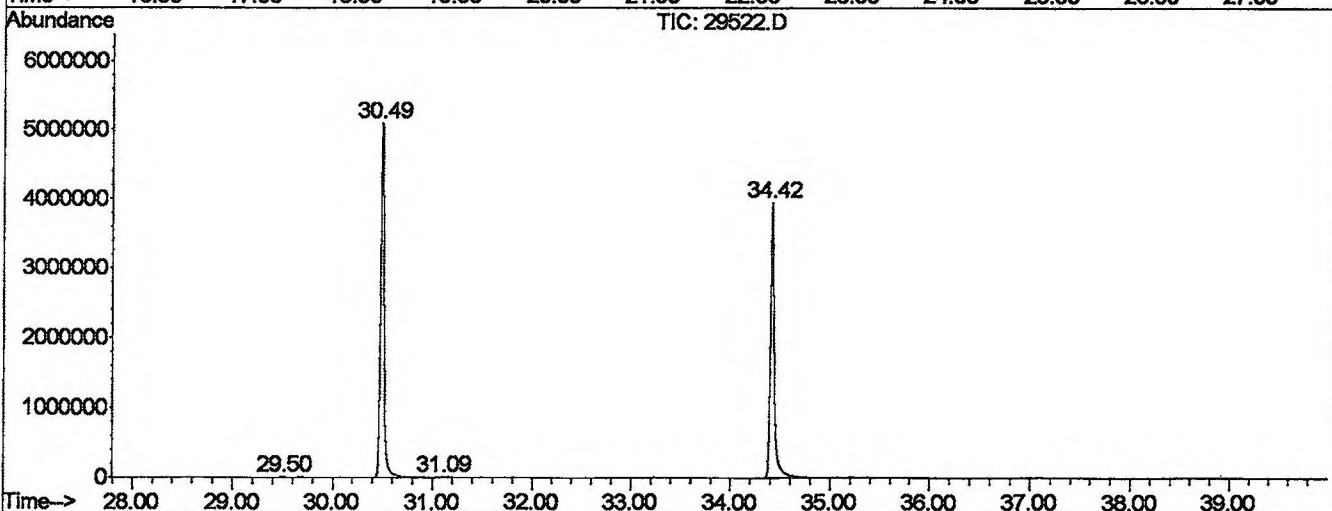
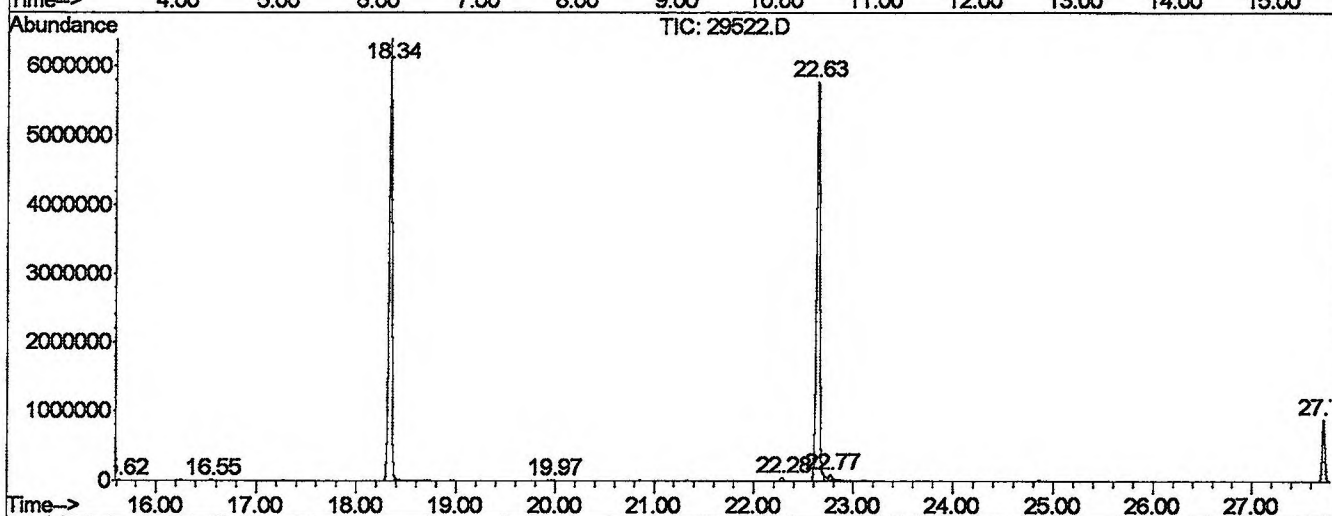
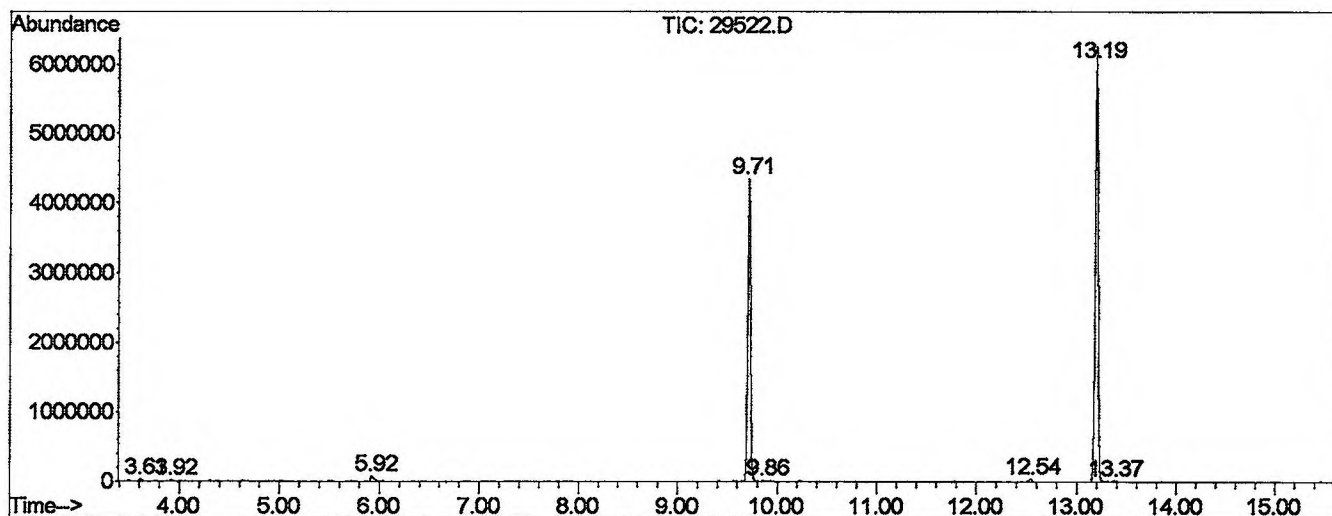
Signal : TIC

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.613	19	21	25	rBV	33342	53348	0.39%	0.073%
2	3.922	45	48	53	rBV	17601	29463	0.22%	0.040%
3	5.920	221	223	239	rBV	78979	195775	1.44%	0.268%
4	9.710	551	555	561	rVV	4343769	8487382	62.58%	11.615%
5	9.858	561	568	575	rVB	14104	30865	0.23%	0.042%
6	12.541	799	803	807	rVV	41045	67119	0.49%	0.092%
7	13.192	855	860	865	rBV	6260394	11976265	88.31%	16.390%
8	13.375	873	876	881	rVB2	13983	26241	0.19%	0.036%
9	15.624	1067	1073	1077	rBV2	14226	27364	0.20%	0.037%
10	16.549	1151	1154	1159	rVB2	15266	31444	0.23%	0.043%
11	18.341	1305	1311	1315	rBV	6390280	12843050	94.70%	17.576%
12	19.974	1451	1454	1465	rVV3	10575	24655	0.18%	0.034%
13	22.280	1653	1656	1661	rVV2	50316	98020	0.72%	0.134%
14	22.634	1681	1687	1693	rBV2	5769957	13562220	100.00%	18.560%
15	22.771	1697	1699	1705	rVB	79935	161468	1.19%	0.221%
16	27.726	2129	2133	2139	rBV	893435	1596218	11.77%	2.184%
17	29.496	2283	2288	2295	rBV3	15868	38889	0.29%	0.053%
18	30.489	2369	2375	2411	rBV	5099453	13117119	96.72%	17.951%
19	31.094	2425	2428	2435	rVV	13029	36925	0.27%	0.051%
20	34.416	2713	2719	2757	rBV	3953781	10668393	78.66%	14.600%

Sum of corrected areas: 73072223

LSC Report - Integrated Chromatogram

File : C:\MSDCHEM\1\5973N\02FEB01\29522.D
 Operator :
 Acquired : 1 Feb 2002 1:04 pm using AcqMethod SCAN508
 Instrument : Instrumen
 Sample Name: 508 scan
 Misc Info : February 1, 2002
 Vial Number: 4
 Quant File :SCAN508.RES (RTE Integrator)



Library Search Compound Report

Data File : C:\MSDCHEM\1\5973N\02FEB01\29522.D
Acq On : 1 Feb 2002 1:04 pm
Sample : 508 scan
Misc : February 1, 2002
MS Integration Params: LSCINT.P

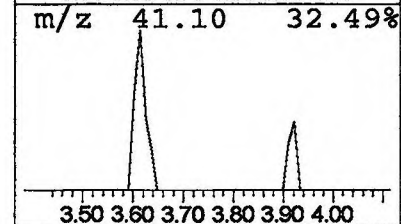
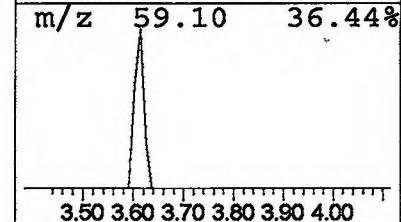
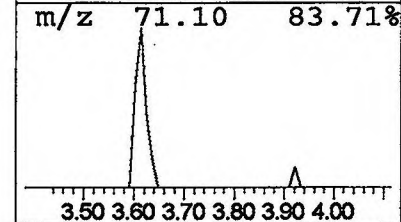
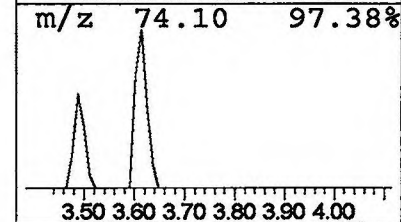
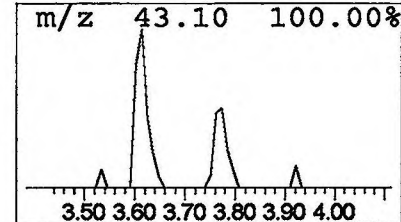
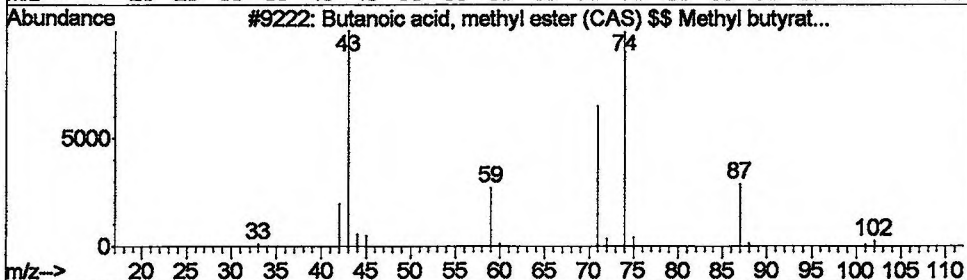
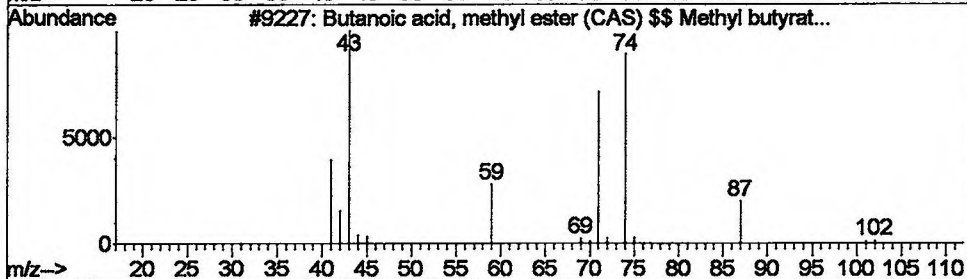
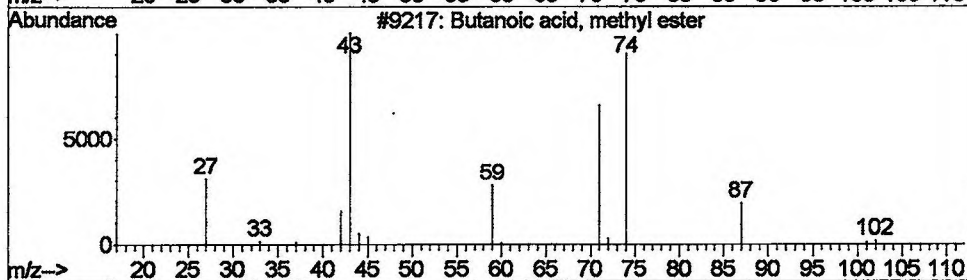
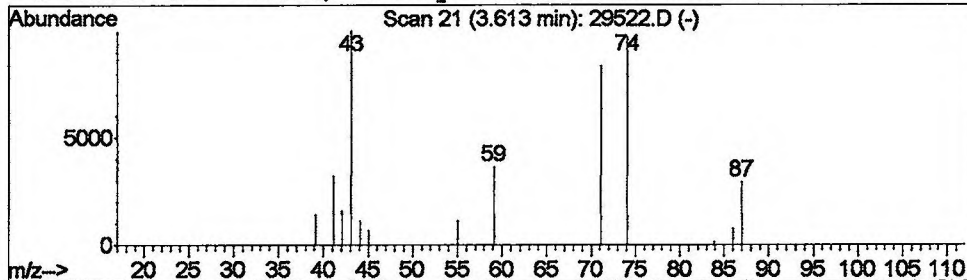
Vial: 4
Operator:
Inst : Instrumen
Multiplr: 1.00

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 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.61	0.13 ug/L	53348	1,4-Dichlorobenzene-d4	9.71

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Butanoic acid, methyl ester	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	78



Library Search Compound Report

Data File : C:\MSDCHEM\1\5973N\02FEB01\29522.D
Acq On : 1 Feb 2002 1:04 pm
Sample : 508 scan
Misc : February 1, 2002
MS Integration Params: LSCINT.P

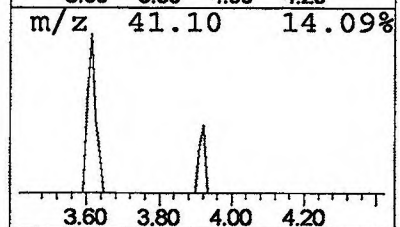
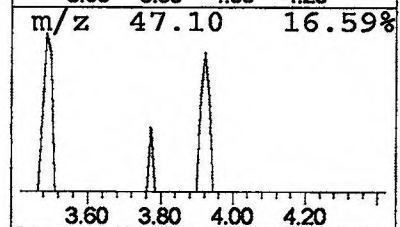
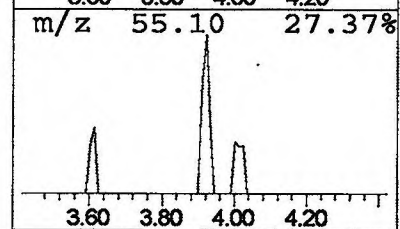
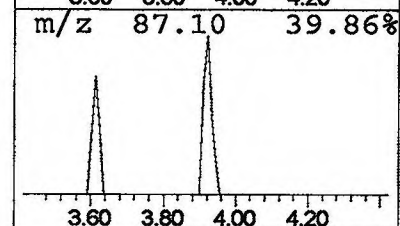
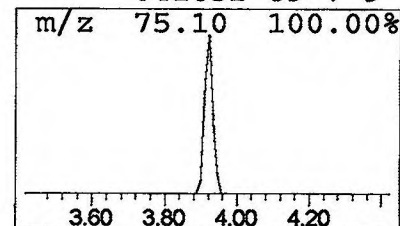
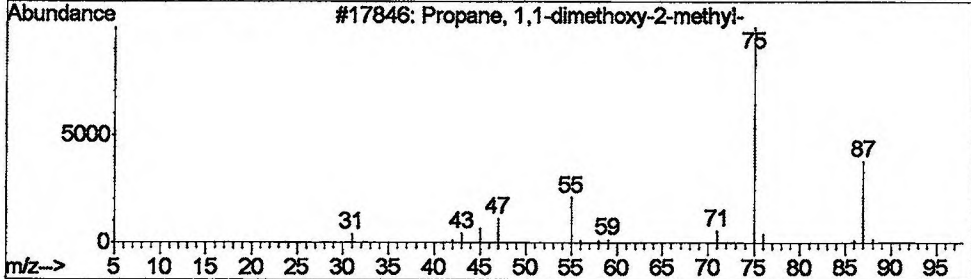
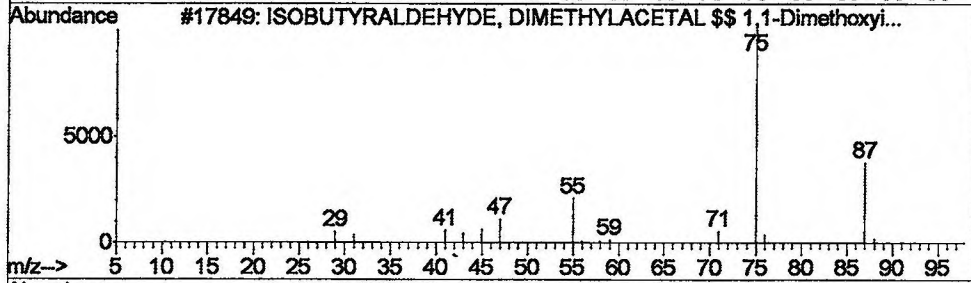
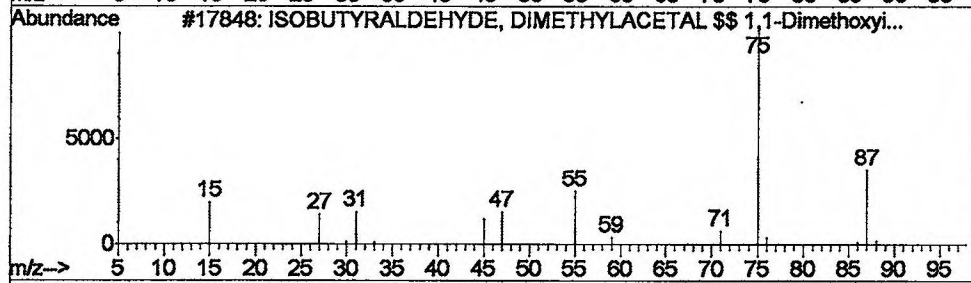
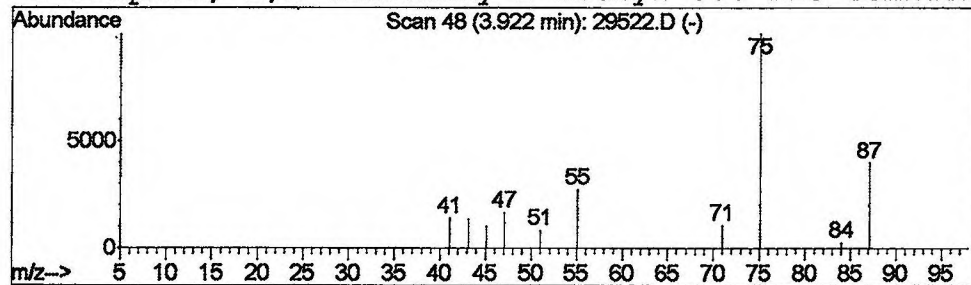
Vial: 4
Operator:
Inst : Instrumen
Multiplr: 1.00

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.92	0.07 ug/L	29463	1,4-Dichlorobenzene-d4	9.71

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			ISOBUTYRALDEHYDE, DIMETHYLACETAL...	118	C6H14O2	000000-00-0	83
2			ISOBUTYRALDEHYDE, DIMETHYLACETAL...	118	C6H14O2	000000-00-0	74
3			Propane, 1,1-dimethoxy-2-methyl-	118	C6H14O2	041632-89-7	74
4			Propane, 1,1-dimethoxy-2-methyl-...	118	C6H14O2	041632-89-7	9



Library Search Compound Report

Data File : C:\MSDCHEM\1\5973N\02FEB01\29522.D
Acq On : 1 Feb 2002 1:04 pm
Sample : 508 scan
Misc : February 1, 2002
MS Integration Params: LSCINT.P

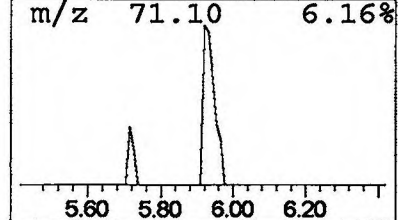
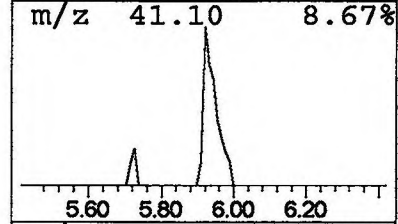
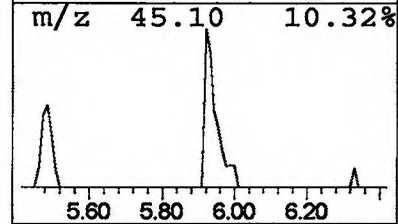
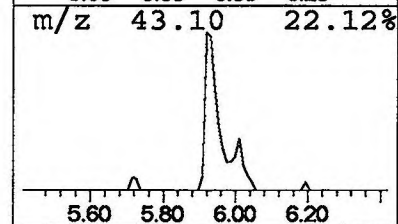
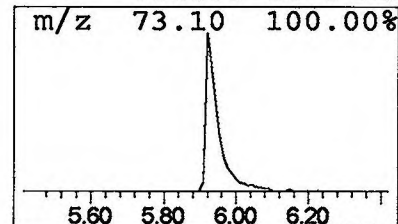
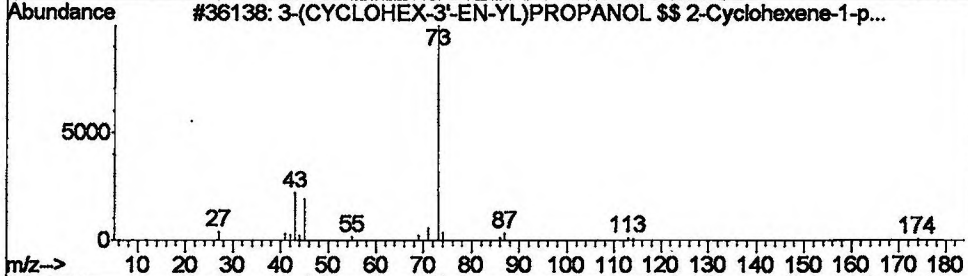
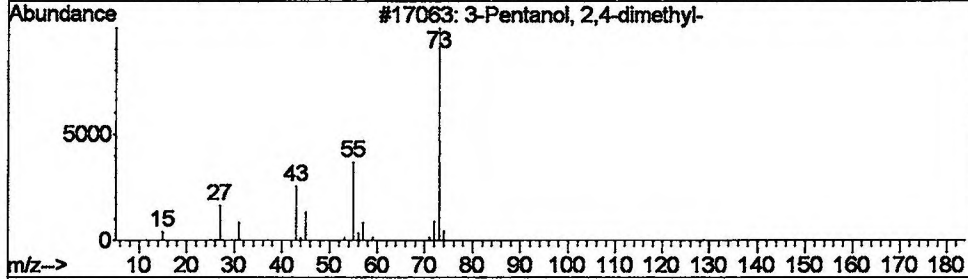
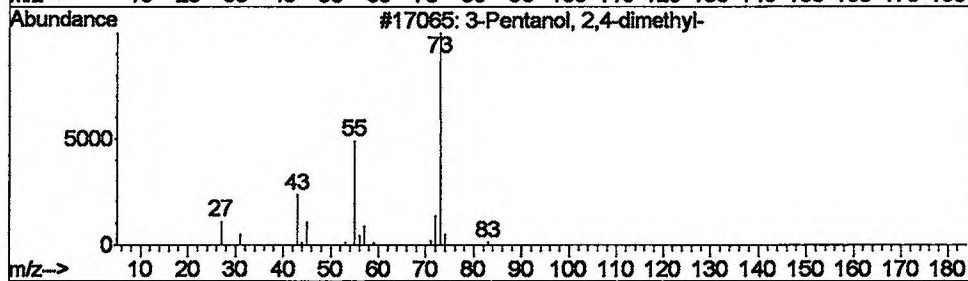
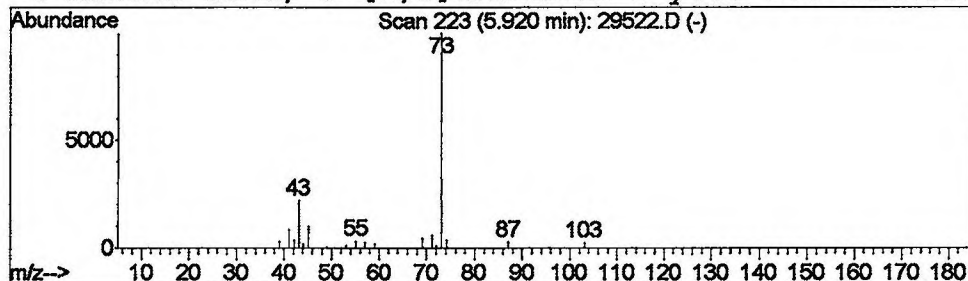
Vial: 4
Operator:
Inst : Instrumen
Multiplr: 1.00

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

Peak Number 3 3-Pentanol, 2,4-dimethyl- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.92	0.46 ug/L	195775	1,4-Dichlorobenzene-d4	9.71

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			3-Pentanol, 2,4-dimethyl-	116	C7H16O	000600-36-2	33
2			3-Pentanol, 2,4-dimethyl-	116	C7H16O	000600-36-2	33
3			3-(CYCLOHEX-3'-EN-YL)PROPANOL \$\$...	174	C8H14O4	015745-87-6	28
4			Acetic acid, 3-[1,3]dioxolan-2-y...	174	C8H14O4	000000-00-0	28



Library Search Compound Report

Data File : C:\MSDCHEM\1\5973N\02FEB01\29522.D
Acq On : 1 Feb 2002 1:04 pm
Sample : 508 scan
Misc : February 1, 2002
MS Integration Params: LSCINT.P

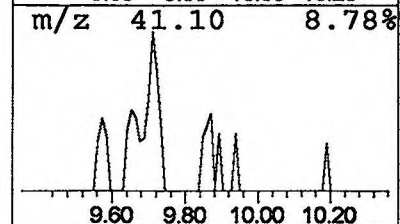
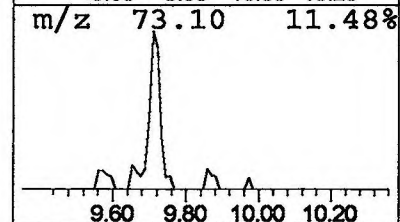
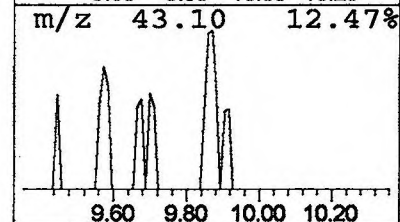
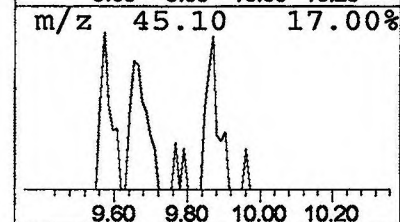
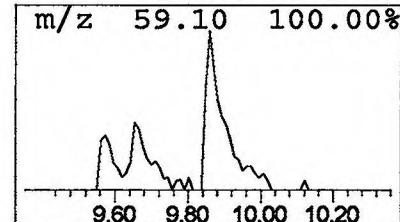
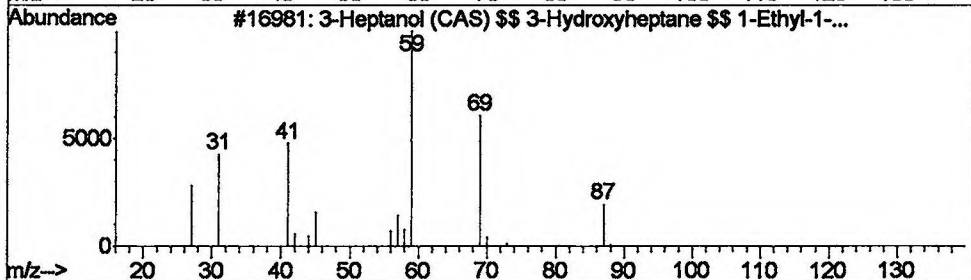
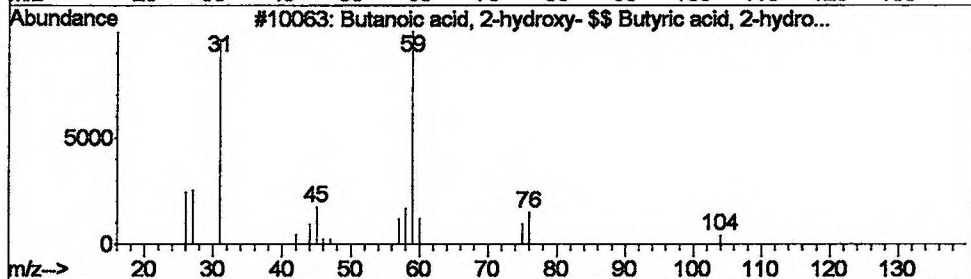
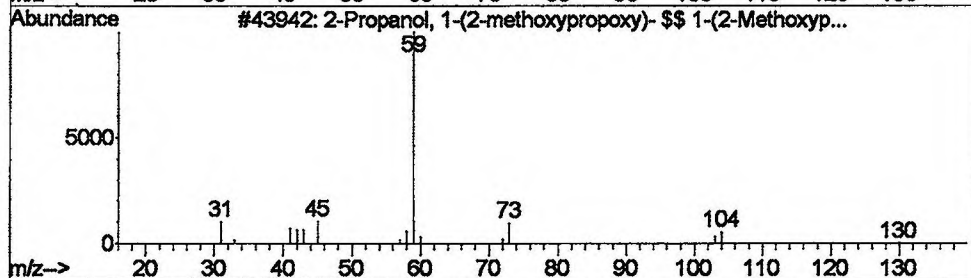
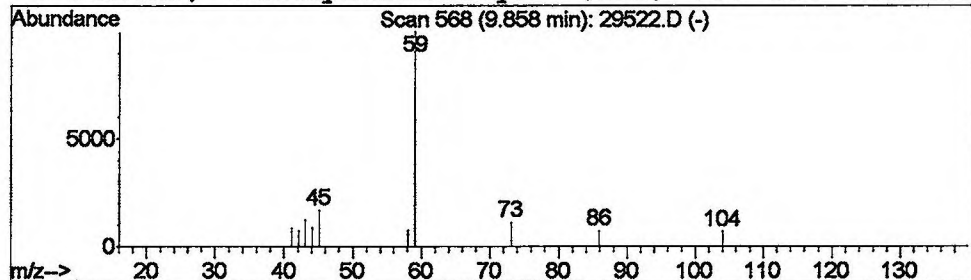
Vial: 4
Operator:
Inst : Instrumen
Multiplr: 1.00

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

Peak Number 4 2-Propanol, 1-(2-methoxypro... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.86	0.07 ug/L	30865	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	74
2			Butanoic acid, 2-hydroxy- \$\$ But...	104	C4H8O3	000565-70-8	39
3			3-Heptanol (CAS) \$\$ 3-Hydroxyhep...	116	C7H16O	000589-82-2	9
4			Silane, ethenyltrimethyl- (CAS) ...	100	C5H12Si	000754-05-2	9



Library Search Compound Report

Data File : C:\MSDCHEM\1\5973N\02FEB01\29522.D
 Acq On : 1 Feb 2002 1:04 pm
 Sample : 508 scan
 Misc : February 1, 2002
 MS Integration Params: LSCINT.P

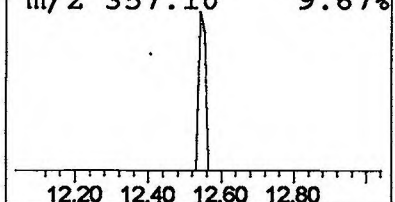
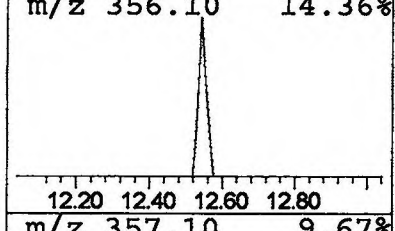
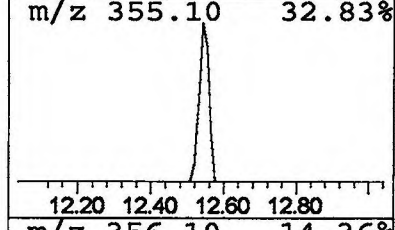
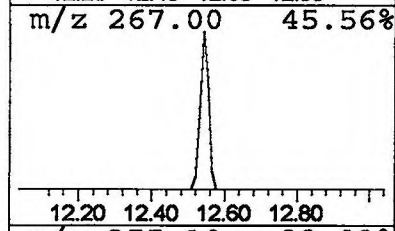
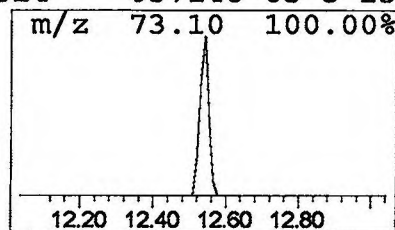
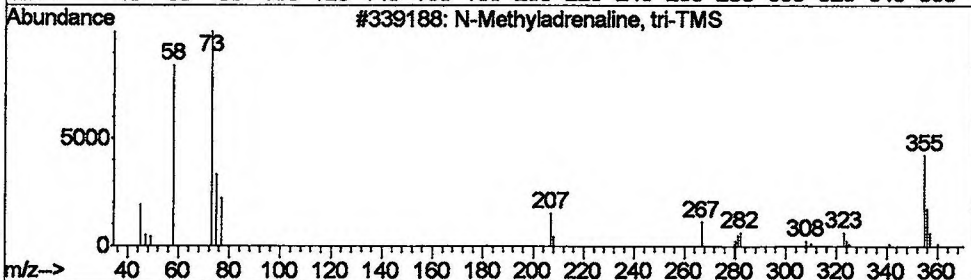
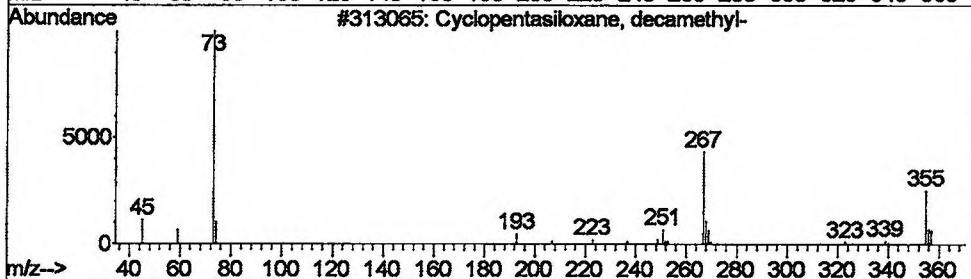
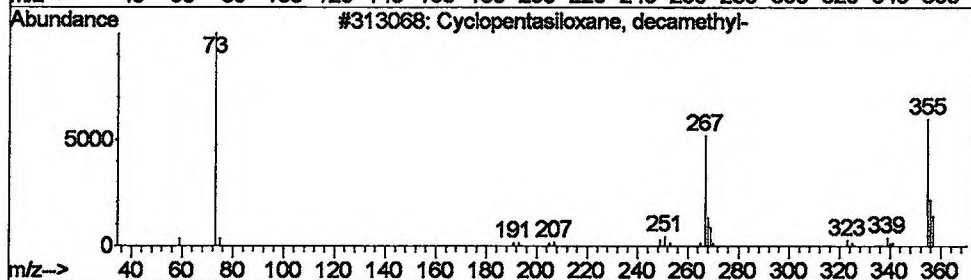
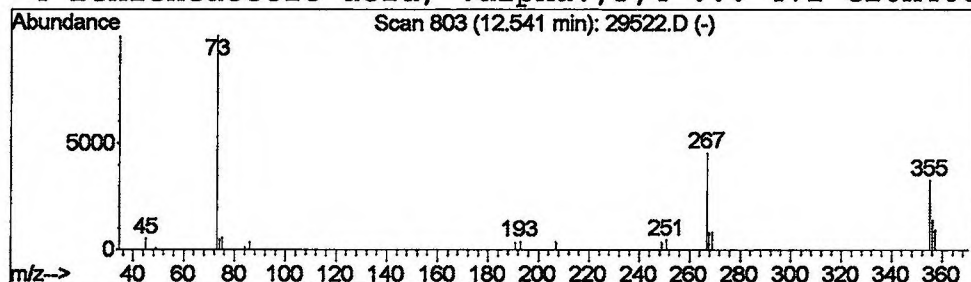
Vial: 4
 Operator:
 Inst : Instrumen
 Multiplr: 1.00

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

 Peak Number 5 Cyclopentasiloxane, decamet... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.54	0.11 ug/L	67119	Naphthalene-d8	13.19

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclopentasiloxane, decamethyl-	370	C10H30O5Si5	000541-02-6	53
2			Cyclopentasiloxane, decamethyl-	370	C10H30O5Si5	000541-02-6	33
3			N-Methyladrenaline, tri-TMS	413	C19H39NO3Si3	000000-00-0	23
4			Benzeneacetic acid, .alpha.,3,4-...	472	C20H40O5Si4	037148-65-5	23



Library Search Compound Report

Data File : C:\MSDCHEM\1\5973N\02FEB01\29522.D
Acq On : 1 Feb 2002 1:04 pm
Sample : 508 scan
Misc : February 1, 2002
MS Integration Params: LSCINT.P

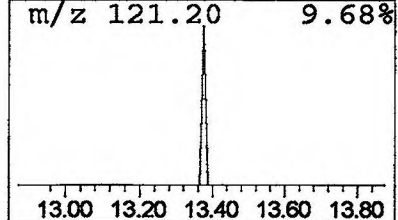
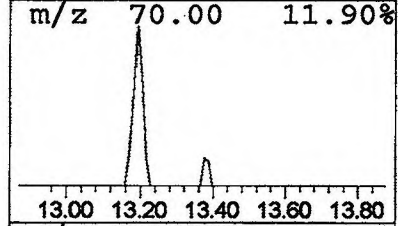
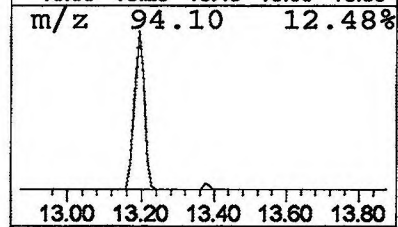
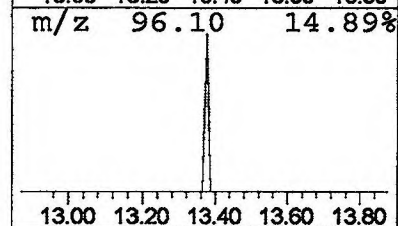
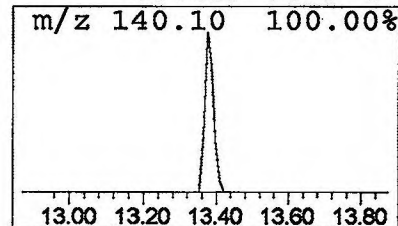
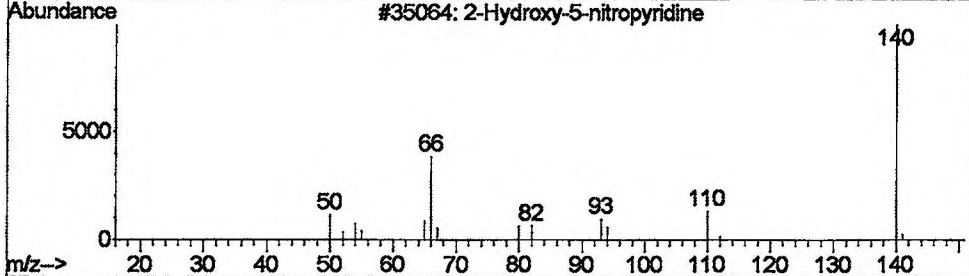
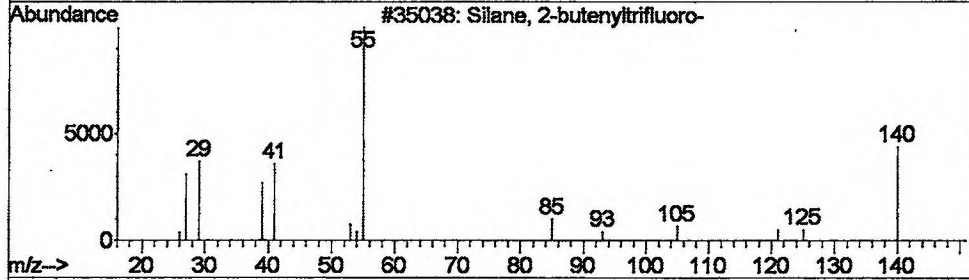
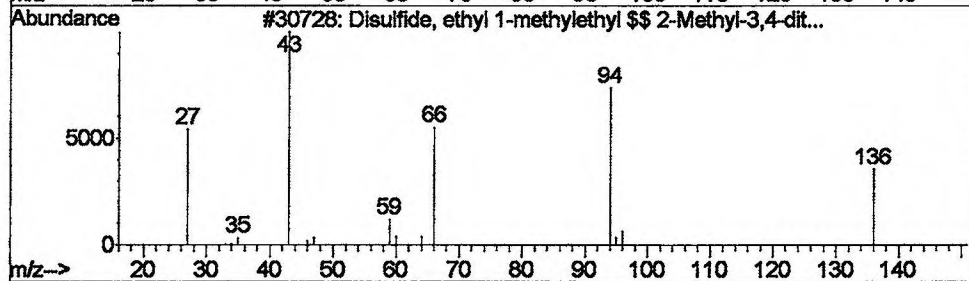
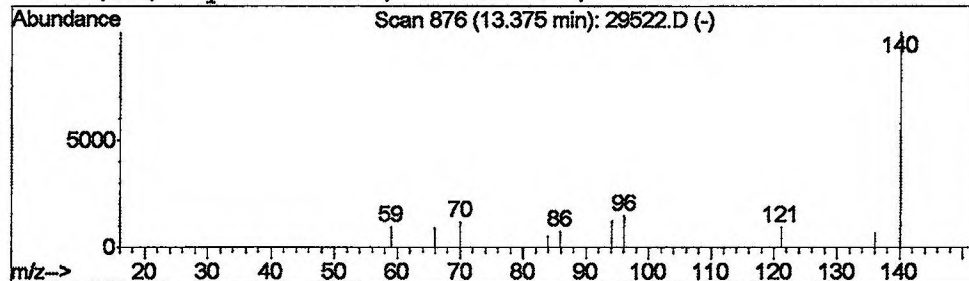
Vial: 4
Operator:
Inst : Instrumen
Multiplr: 1.00

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

Peak Number 6 Disulfide, ethyl 1-methylet... Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.37	0.04 ug/L	26241	Naphthalene-d8	13.19

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Disulfide, ethyl 1-methylethyl \$...	136	C5H12S2	053966-36-2	7
2			Silane, 2-butenyltrifluoro-	140	C4H7F3Si	055712-74-8	4
3			2-Hydroxy-5-nitropyridine	140	C5H4N2O3	005418-51-9	4
4			2(5H)-Pyridinone, 5-diazo-, mono...	121	C5H3N3O	054459-87-9	4



Library Search Compound Report

Data File : C:\MSDCHEM\1\5973N\02FEB01\29522.D
Acq On : 1 Feb 2002 1:04 pm
Sample : 508 scan
Misc : February 1, 2002
MS Integration Params: LSCINT.P

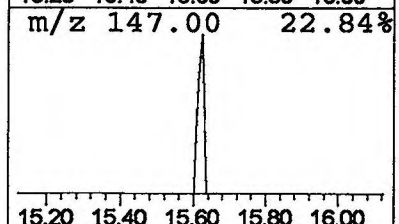
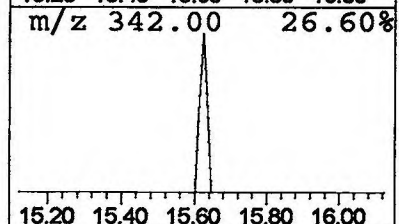
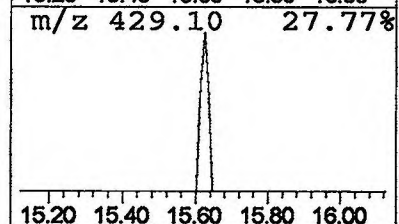
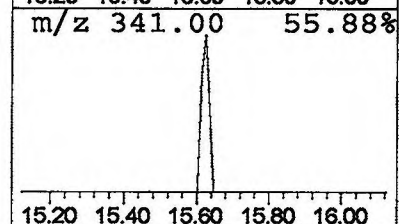
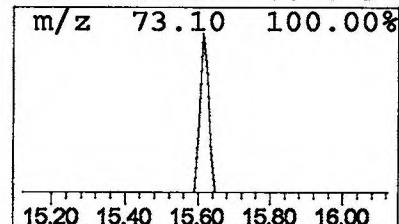
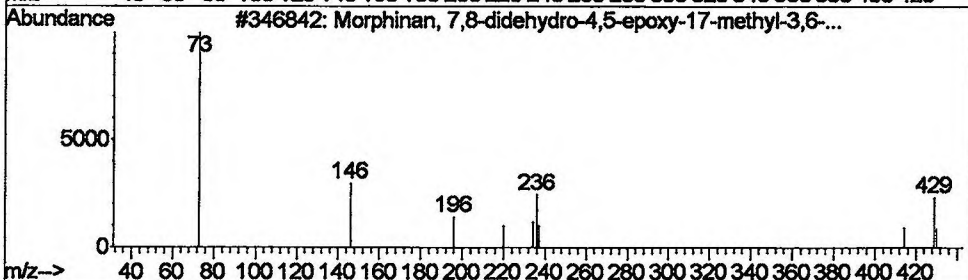
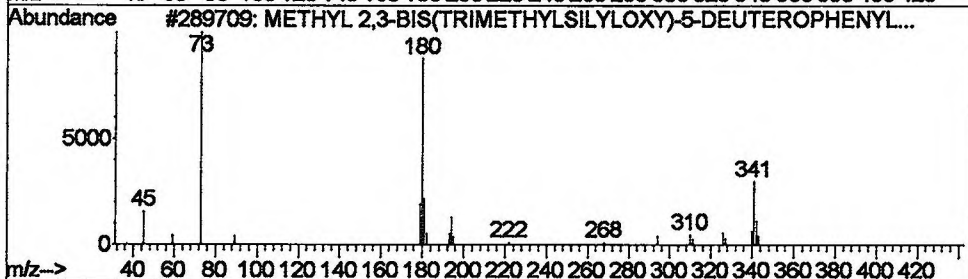
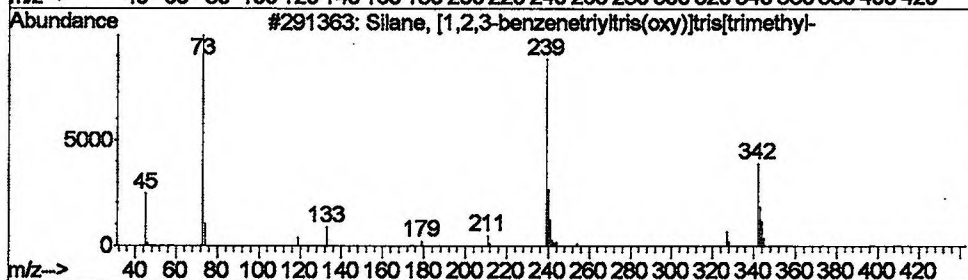
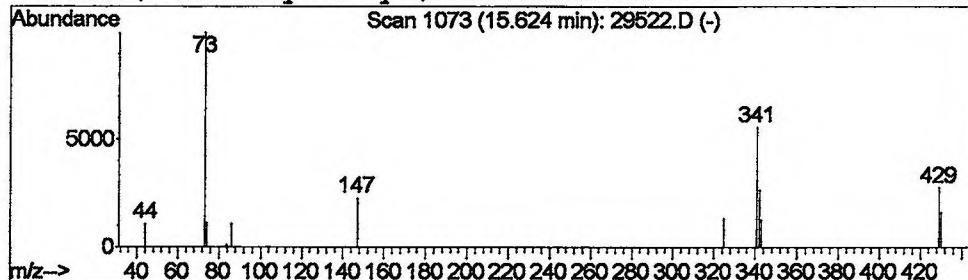
Vial: 4
Operator:
Inst : Instrumen
Multiplr: 1.00

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

Peak Number 7 Silane, [1,2,3-benzenetriyl... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.62	0.05 ug/L	27364	Naphthalene-d8	13.19

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Silane, [1,2,3-benzenetriyltris(...	342	C15H30O3Si3	017864-23-2	10
2			METHYL 2,3-BIS(TRIMETHYLSILOXY...	341	C16H27DO4Si2	000000-00-0	9
3			Morphinan, 7,8-didehydro-4,5-epo...	429	C23H35NO3Si2	055449-66-6	9
4			Bis(trimethylsilyl) derivative o...	429	C23H35NO3Si2	000000-00-0	9



Library Search Compound Report

Data File : C:\MSDCHEM\1\5973N\02FEB01\29522.D
 Acq On : 1 Feb 2002 1:04 pm
 Sample : 508 scan
 Misc : February 1, 2002
 MS Integration Params: LSCINT.P

Vial: 4
 Operator:
 Inst : Instrumen
 Multiplr: 1.00

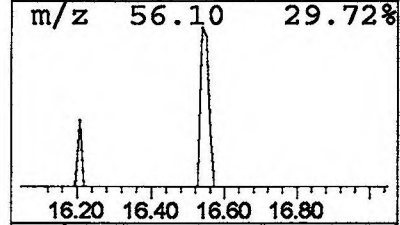
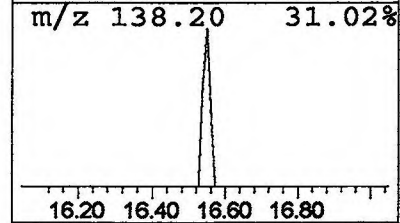
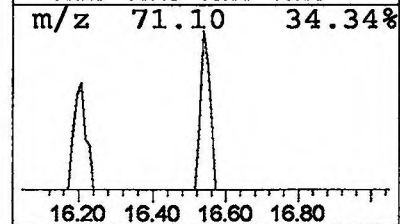
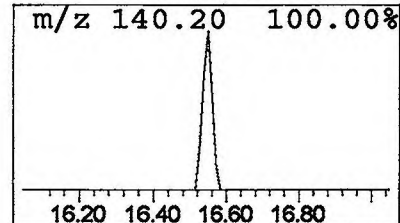
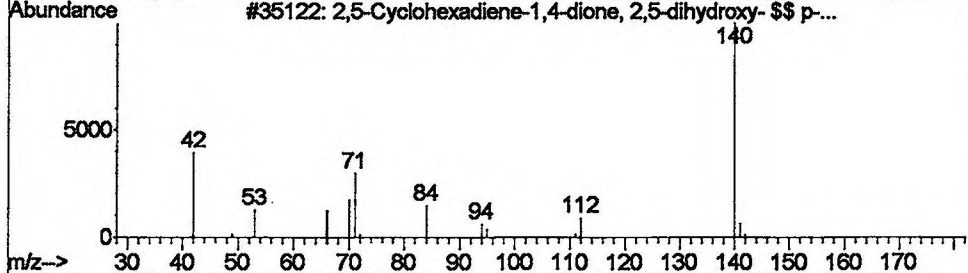
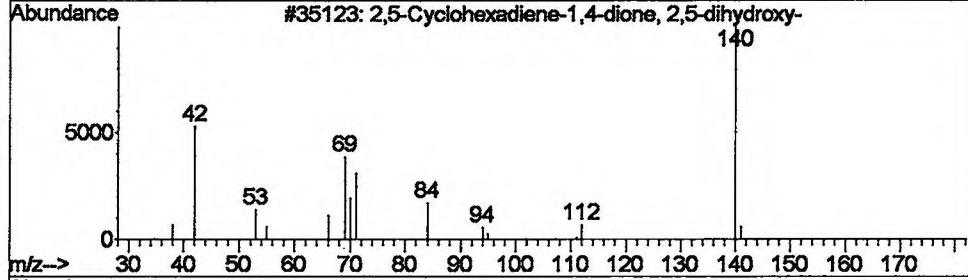
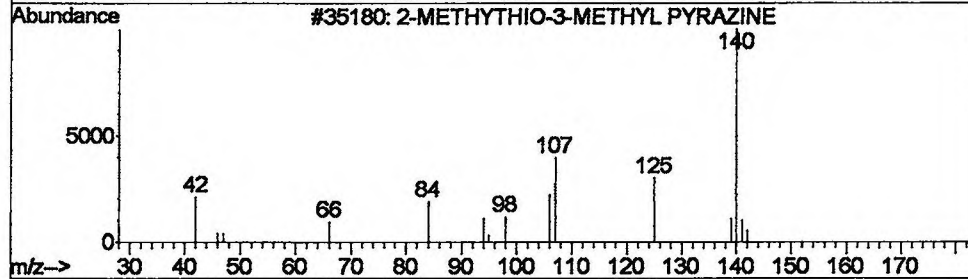
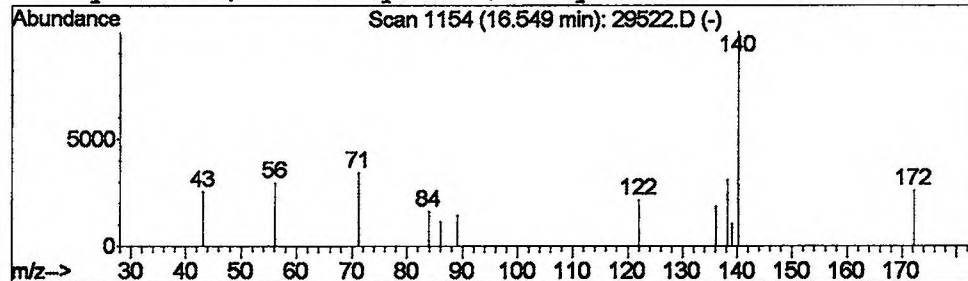
Quant Method : C:\MSDCHEM\1...\SCAN508.M (RTE Integrator)
 Title : 508 scan method
 Library : C:\DATABASE\WILEY7N.L

 Peak Number 8 2-METHYTHIO-3-METHYL PYRAZINE Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.55	0.05 ug/L	31444	Acenaphthene-d10	18.34

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-METHYTHIO-3-METHYL PYRAZINE	140	C6H8N2S	002882-20-4	9
2		2,5-Cyclohexadiene-1,4-dione, 2,...	140	C6H4O4	000615-94-1	9
3		2,5-Cyclohexadiene-1,4-dione, 2,...	140	C6H4O4	000615-94-1	9
4		Pyrazine, 2-methyl-3-(methylthio...	140	C6H8N2S	002882-20-4	9

NO GOOD MATCH



Library Search Compound Report

Data File : C:\MSDCHEM\1\5973N\02FEB01\29522.D
Acq On : 1 Feb 2002 1:04 pm
Sample : 508 scan
Misc : February 1, 2002
MS Integration Params: LSCINT.P

Vial: 4
Operator:
Inst : Instrumen
Multiplr: 1.00

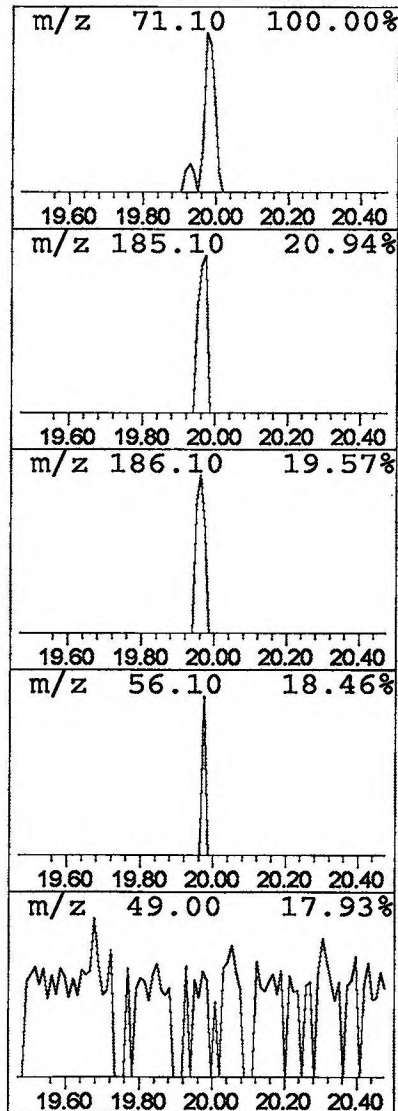
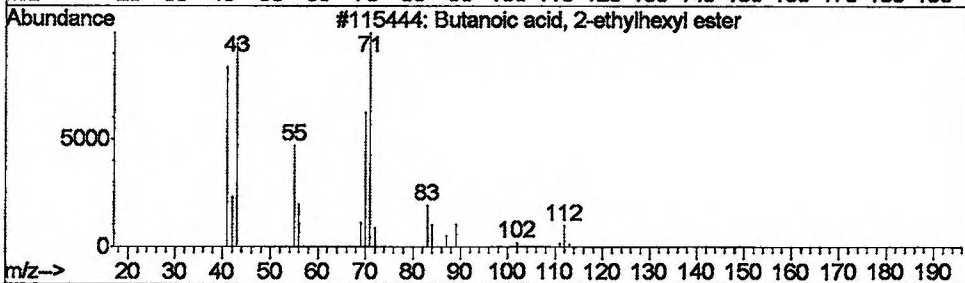
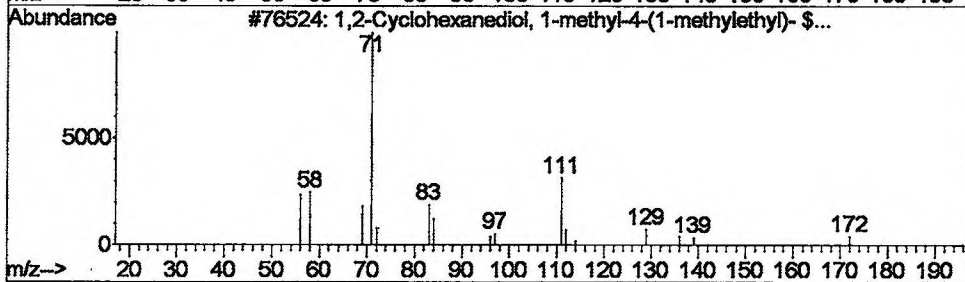
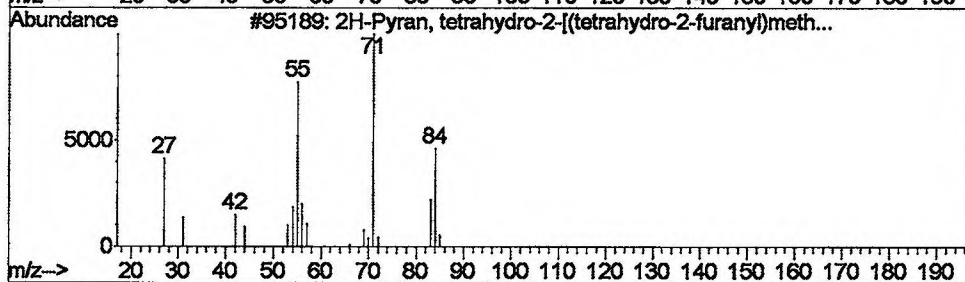
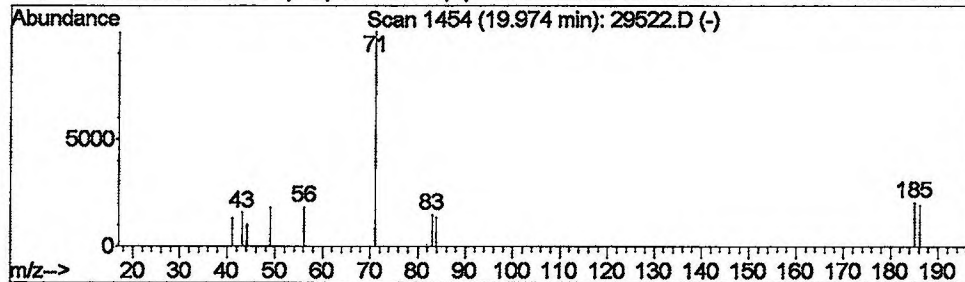
Quant Method : C:\MSDCHEM\1...\SCAN508.M (RTE Integrator)
Title : 508 scan method
Library : C:\DATABASE\WILEY7N.L

Peak Number 9 2H-Pyran, tetrahydro-2-[(te... Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.97	0.04 ug/L	24655	Acenaphthene-d10	18.34

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2H-Pyran, tetrahydro-2-[(tetrahy...	186	C10H18O3	000710-14-5	9
2			1,2-Cyclohexanediol, 1-methyl-4-...	172	C10H20O2	033669-76-0	9
3			Butanoic acid, 2-ethylhexyl ester	200	C12H24O2	000000-00-0	9
4			Imidazole-2,4,5-d3 \$\$ 1H-Imidazo...	71	C3HD3N2	006745-43-3	4

ND
Good
match



Library Search Compound Report

Data File : C:\MSDCHEM\1\5973N\02FEB01\29522.D

Acq On : 1 Feb 2002 1:04 pm

Sample : 508 scan

Misc : February 1, 2002

MS Integration Params: LSCINT.P

Vial: 4

Operator:

Inst : Instrumen

Multiplr: 1.00

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

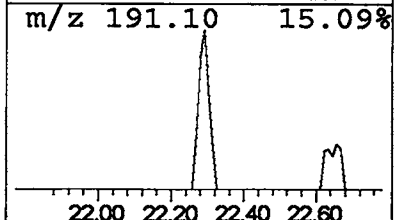
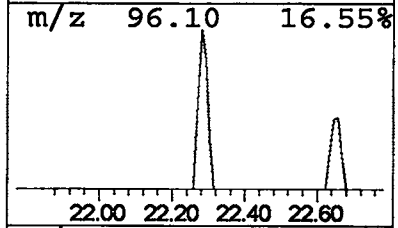
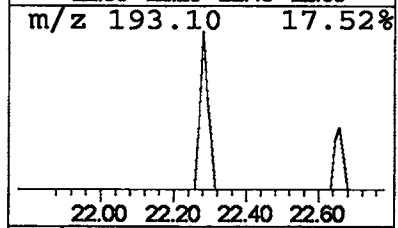
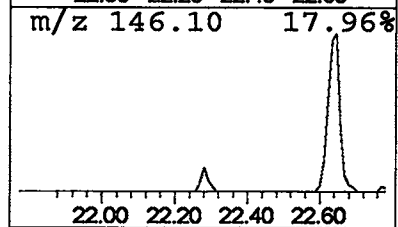
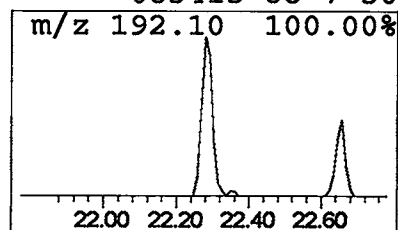
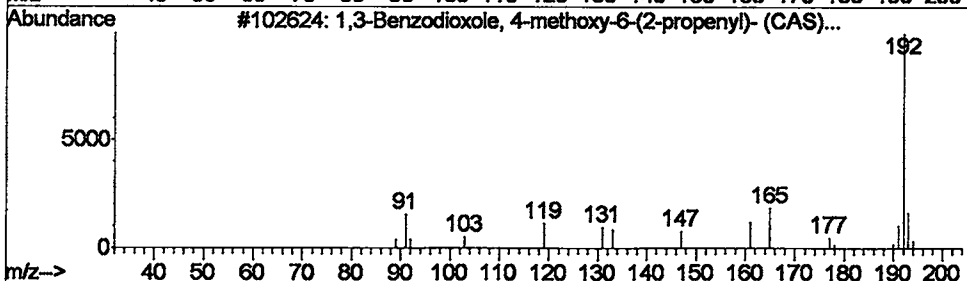
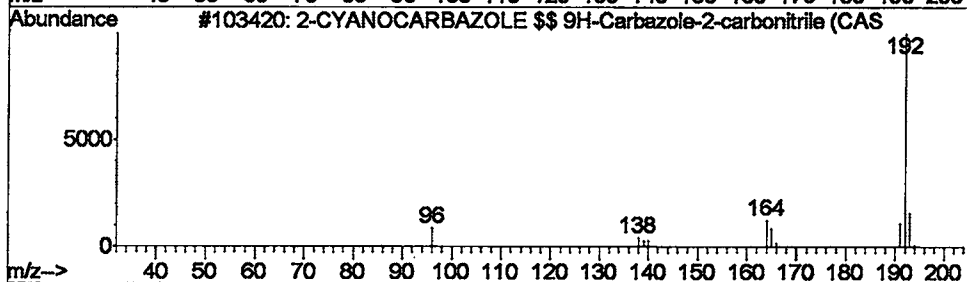
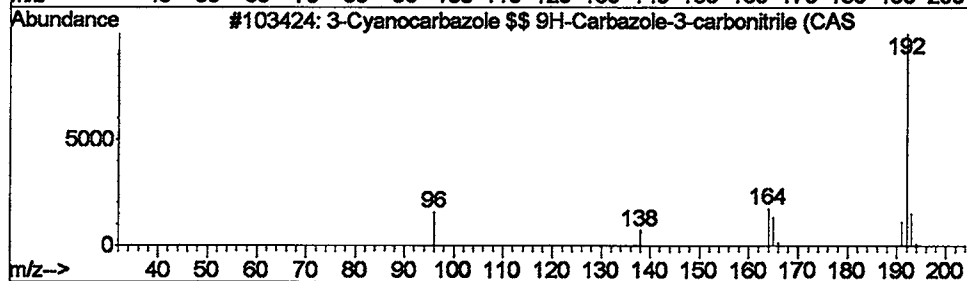
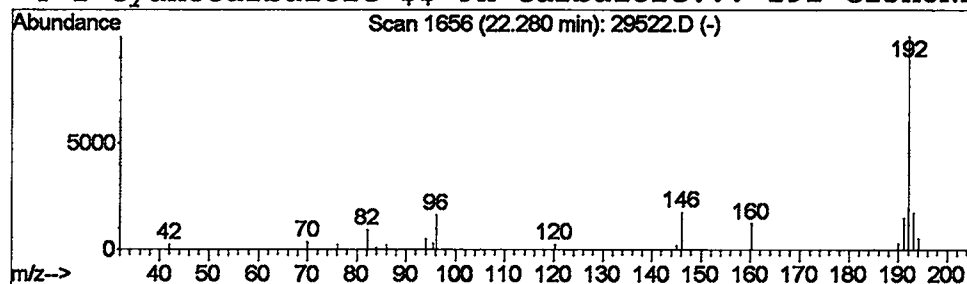
Title : 508 scan method

Library : C:\DATABASE\WILEY7N.L

Peak Number 10 3-Cyanocarbazole \$\$ 9H-Carb... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.28	0.14 ug/L	98020	Phenanthrene-d10	22.63

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		3-Cyanocarbazole \$\$ 9H-Carbazole...	192	C13H8N2	057102-93-9	72
2		2-CYANOCARBAZOLE \$\$ 9H-Carbazole...	192	C13H8N2	057955-18-7	72
3		1,3-Benzodioxole, 4-methoxy-6-(2...	192	C11H12O3	000607-91-0	53
4		1-Cyanocarbazole \$\$ 9H-Carbazole...	192	C13H8N2	083415-88-7	50



Library Search Compound Report

Data File : C:\MSDCHEM\1\5973N\02FEB01\29522.D
Acq On : 1 Feb 2002 1:04 pm
Sample : 508 scan
Misc : February 1, 2002
MS Integration Params: LSCINT.P

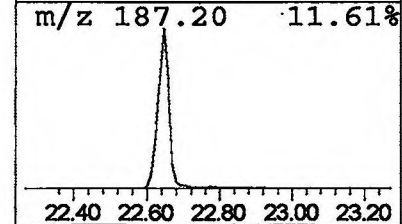
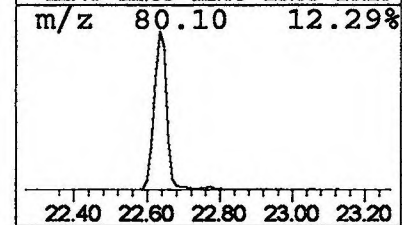
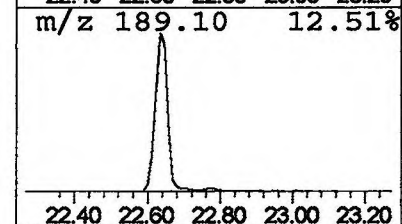
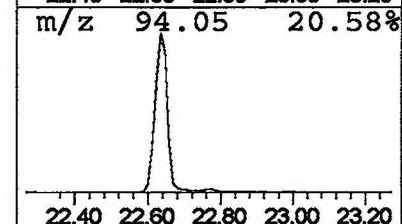
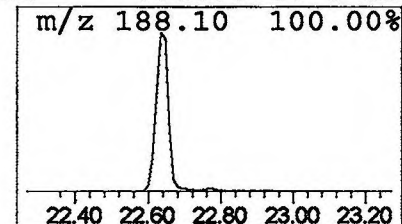
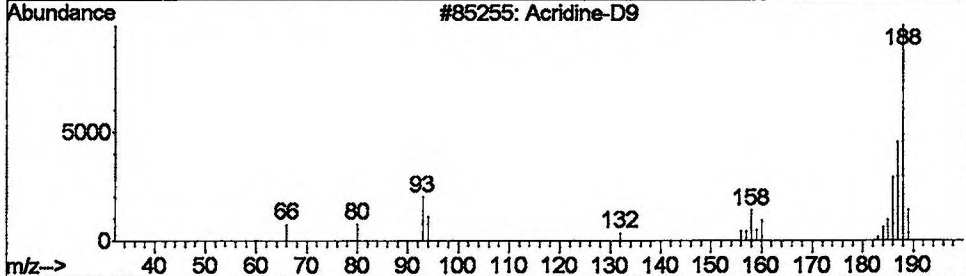
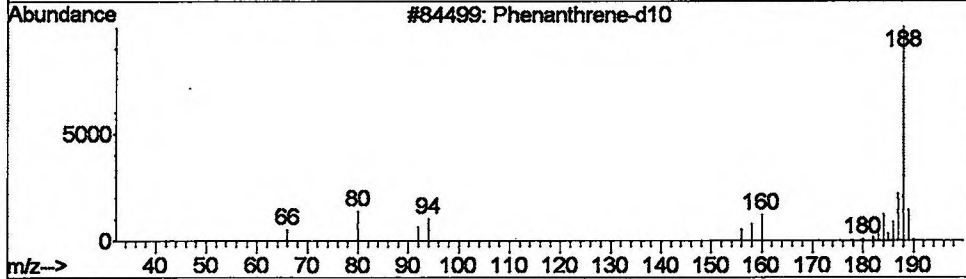
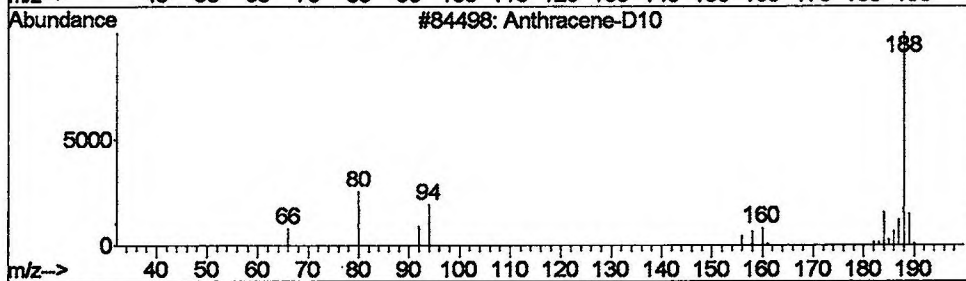
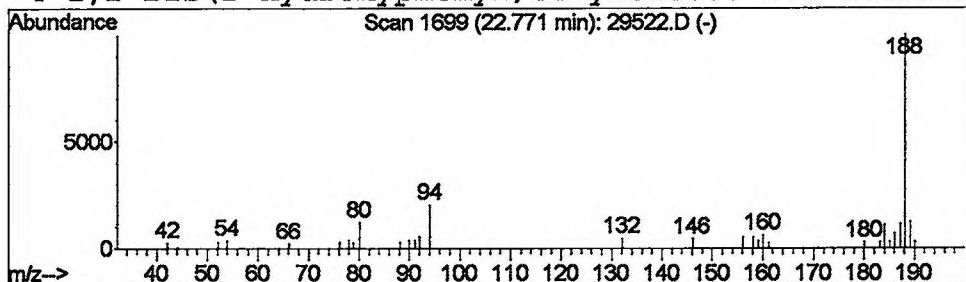
Vial: 4
Operator:
Inst : Instrumen
Multiplr: 1.00

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 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.77	0.24 ug/L	161468	Phenanthrene-d10	22.63

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Anthracene-D10	188	C14D10	000000-00-0	80
2			Phenanthrene-d10	188	C14D10	001517-22-2	68
3			Acridine-D9	188	C13D9N	000000-00-0	64
4			1,2-Bis(2-hydroxyphenyl)ethylene...	376	C24H28N2O2	000000-00-0	59



Library Search Compound Report

Data File : C:\MSDCHEM\1\5973N\02FEB01\29522.D
 Acq On : 1 Feb 2002 1:04 pm
 Sample : 508 scan
 Misc : February 1, 2002
 MS Integration Params: LSCINT.P

Vial: 4
 Operator:
 Inst : Instrumen
 Multiplr: 1.00

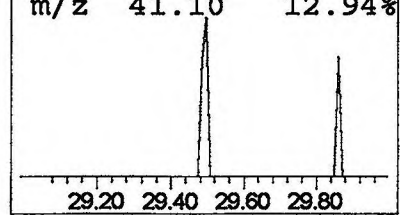
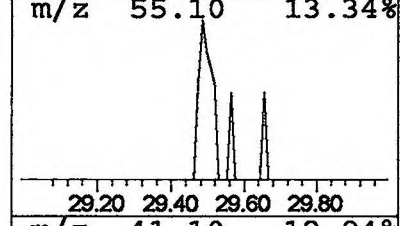
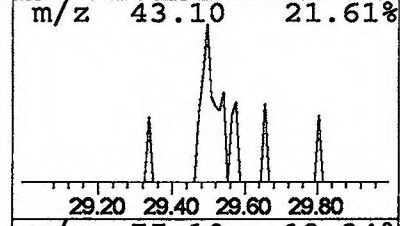
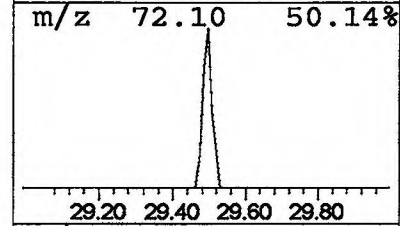
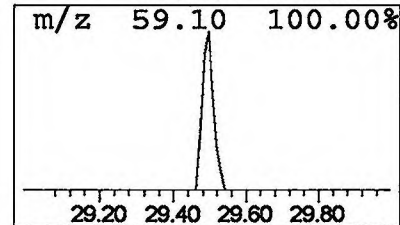
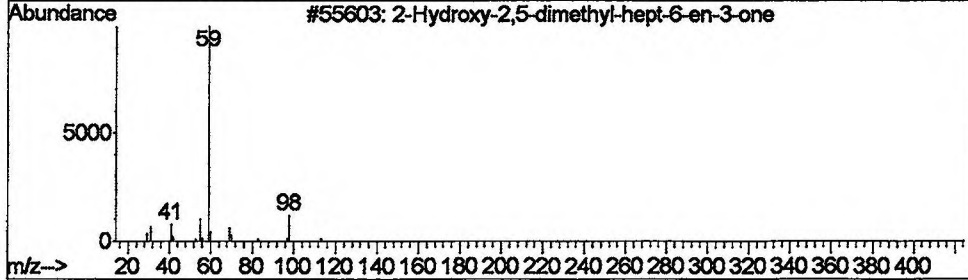
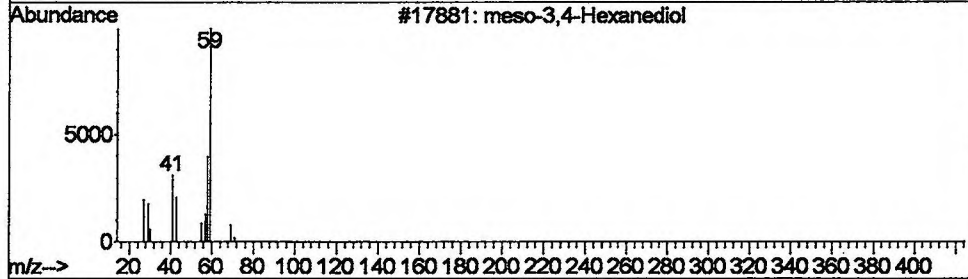
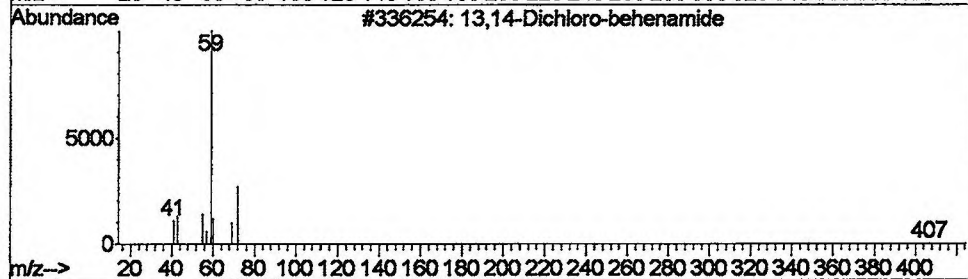
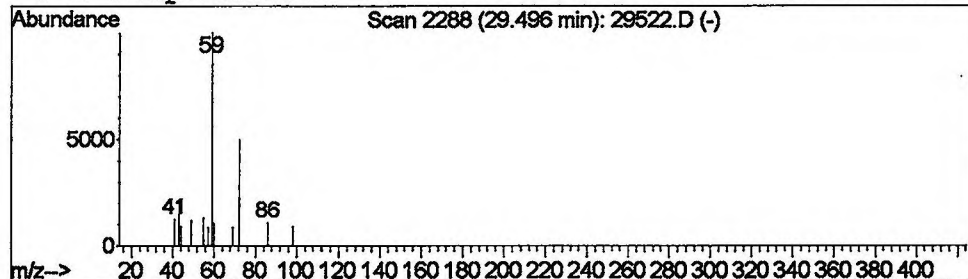
Quant Method : C:\MSDCHEM\1...\SCAN508.M (RTE Integrator)
 Title : 508 scan method
 Library : C:\DATABASE\WILEY7N.L

 Peak Number 12 13,14-Dichloro-behenamide Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
29.50	0.06 ug/L	38889	Chrysene-d12	30.50

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			13,14-Dichloro-behenamide	407	C22H43Cl2NO	000000-00-0	33
2			meso-3,4-Hexanediol	118	C6H14O2	022520-39-4	9
3			2-Hydroxy-2,5-dimethyl-hept-6-en...	156	C9H16O2	000000-00-0	9
4			1-Propanamine	59	C3H9N	000107-10-8	7

*NO
good
match*



Tentatively Identified Compound (LSC) summary

Operator ID: Date Acquired: 1 Feb 2002 1:04 pm
 Data File: C:\MSDCHEM\1\5973N\02FEB01\29522.D
 Name: 508 scan
 Misc: February 1, 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...	3.61	0.1	ug/L	53348	1	9.71	8487380	20.0
ISOBUTYRALDEHYDE...	3.92	0.1	ug/L	29463	1	9.71	8487380	20.0
3-Pentanol, 2,4-d...	5.92	0.5	ug/L	195775	1	9.71	8487380	20.0
2-Propanol, 1-(2-...	9.86	0.1	ug/L	30865	1	9.71	8487380	20.0
Cyclopentasiloxan...	12.54	0.1	ug/L	67119	2	13.19	11976300	20.0
Disulfide, ethyl ...	13.37	0.0	ug/L	26241	2	13.19	11976300	20.0
Silane, [1,2,3-be...	15.62	0.0	ug/L	27364	2	13.19	11976300	20.0
2-METHYTHIO-3-MET...	16.55	0.0	ug/L	31444	3	18.34	12843100	20.0
2H-Pyran, tetrahy...	19.97	0.0	ug/L	24655	3	18.34	12843100	20.0
3-Cyanocarbazole ...	22.28	0.1	ug/L	98020	4	22.63	13562200	20.0
Anthracene-D10	22.77	0.2	ug/L	161468	4	22.63	13562200	20.0
13,14-Dichloro-be...	29.50	0.1	ug/L	38889	5	30.50	13117100	20.0