

User: Vinci, David L
Westchester Dept. of Labs & Research
Date: 01/29/2002 Time: 12:01:14

Sample: AE00699
Analysis: \$GCMS GCMS Scan For Unknowns
Units: ug
Started: 1/29/02 12:00 Ended: 1/29/02 12:00 Analyst: DLV
Page: 1

	Component Name	Viol	Result	Qualifier	MDL
1	Other unknown compounds		.		
2	Surr_2,4-DDD		103		5.0

Comments for Sample AE00699 - Analysis \$GCMS

Westchester Dept. of Labs & Research

User: Vinci, David L

Date: 01-29-2002 Time: 12:01:35

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

The recoveries for 15 of the 43 compounds in the LCS Standard were outside of their acceptance ranges.

Quantitation Report (Not Reviewed)

Data File : C:\MSDCHEM\1\5973N\02JAN16\00699.D Vial: 4
 Acq On : 16 Jan 2002 1:20 pm Operator:
 Sample : 508scan Inst : Instrumen
 Misc : January 16, 2002 Multiplr: 1.00
 MS Integration Params: SPLIT.P
 Quant Time: Jan 16 14:05:05 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	1122891	20.00	ug/L	0.00
10) Naphthalene-d8	13.19	136	4710465	20.00	ug/L	-0.02
17) Acenaphthene-d10	18.34	164	2474159	20.00	ug/L	-0.01
29) Phenanthrene-d10	22.63	188	4146871	20.00	ug/L	-0.03
38) Chrysene-d12	30.49	240	3630816	20.00	ug/L	-0.02
44) Perylene-d12	34.42	264	2619126	20.00	ug/L	0.00

System Monitoring Compounds
 37) 2,4-DDD 27.73 235 338093 5.16 ug/L 0.00
 Spiked Amount 5.000 Recovery = 103.20%

Target Compounds					Qvalue
2) n-Nitrosodimethylamine	3.60	74	10907	0.23 ug/L #	10
20) Dimethylphthalate	18.34	163	404818	2.48 ug/L #	55
21) 2,6-Dinitrotoluene	18.34	165	312420	9.80 ug/L #	17
35) Di-n-butylphthalate	24.85	149	10526	0.04 ug/L #	76
41) Benzo(a)anthracene	30.49	228	8058	0.04 ug/L #	49
42) Bis(2-ethylhexyl)phthalate	31.11	149	5187	0.58 ug/L #	44
43) Chrysene	30.49	228	8058	0.04 ug/L #	46
48) Benzo(a)pyrene	34.42	252	8272	0.05 ug/L #	1

(#) = qualifier out of range (m) = manual integration
 00699.D SCAN508.M Wed Jan 16 14:05:06 2002

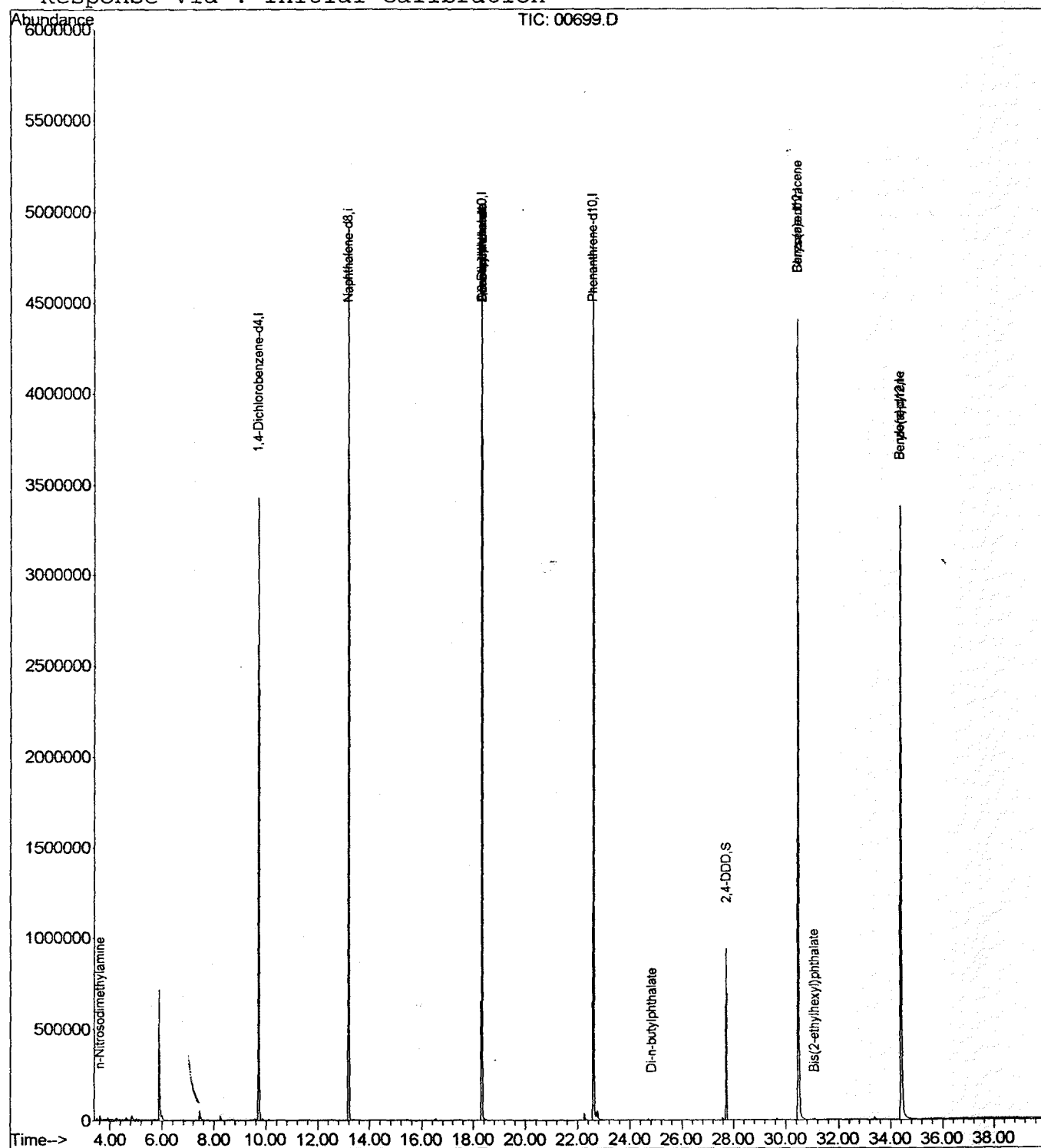
Quantitation Report (Not Reviewed)

Data File : C:\MSDCHEM\1\5973N\02JAN16\00699.D
 Acq On : 16 Jan 2002 1:20 pm
 Sample : 508scan
 Misc : January 16, 2002
 MS Integration Params: SPLIT.P
 Quant Time: Jan 16 14:05 2002

Vial: 4
 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



00699.D SCAN508.M

Wed Jan 16 14:05:07 2002

Page 2

LSC Area Percent Report

Data File : C:\MSDCHEM\1\5973N\02JAN16\00699.D
 Acq On : 16 Jan 2002 1:20 pm
 Sample : 508scan
 Misc : January 16, 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

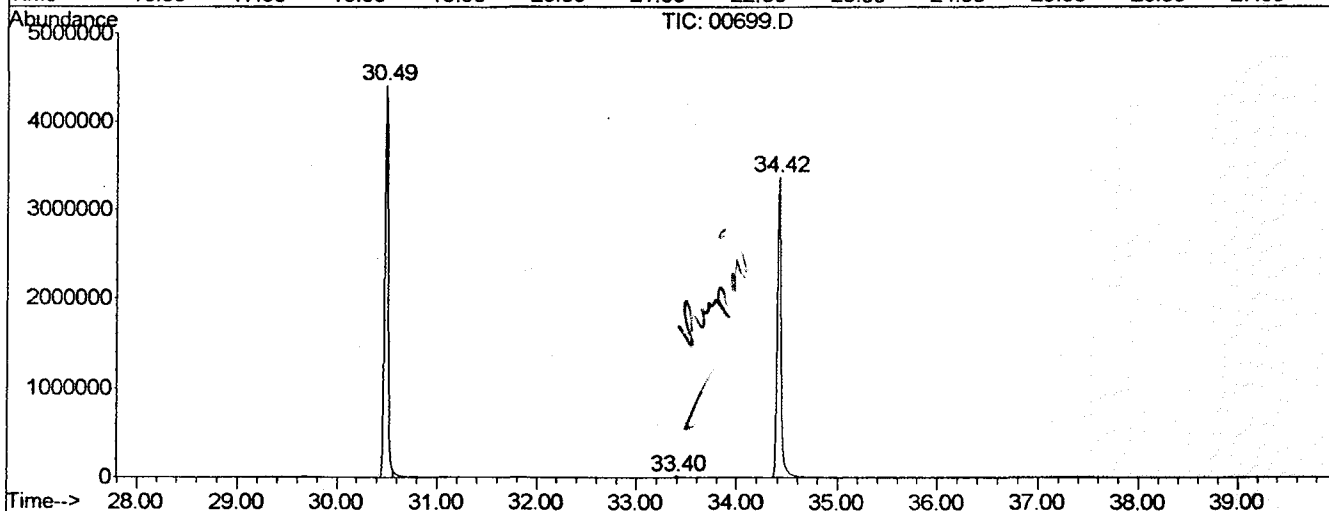
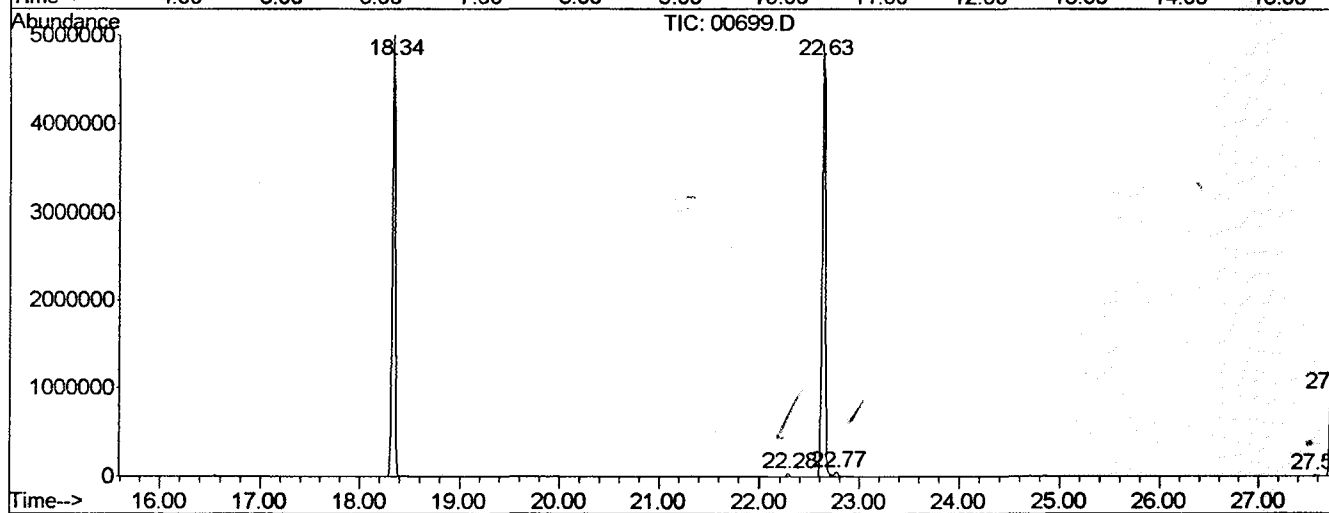
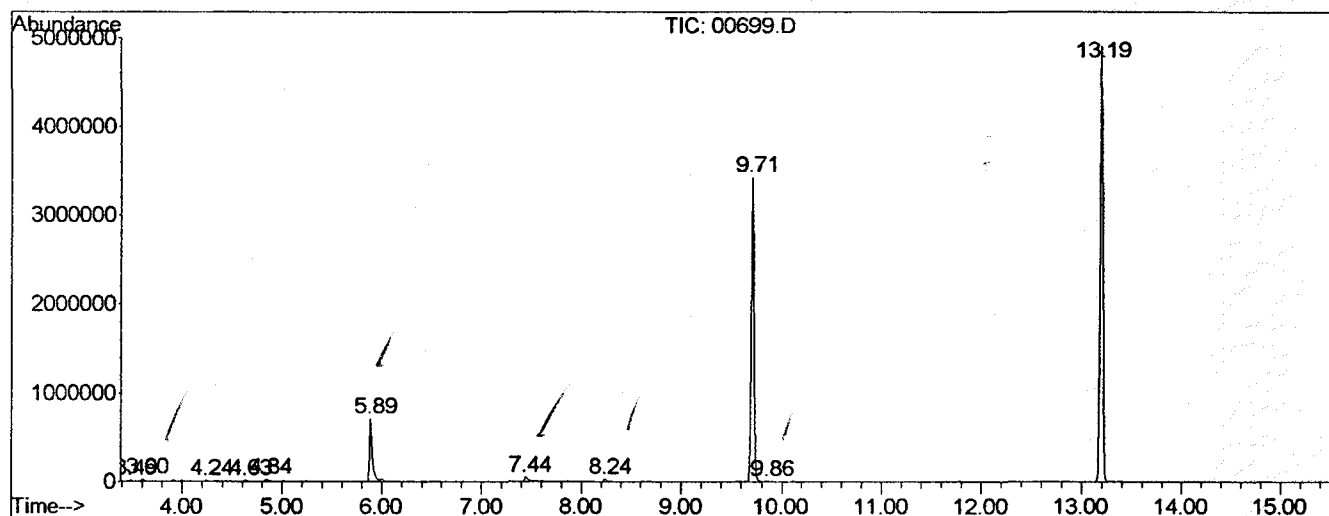
Signal : TIC

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.488	5	10	17	rBV	17039	33887	0.31%	0.057%
2	3.602	17	20	25	rVV	27420	44714	0.41%	0.075%
3	4.241	73	76	83	rVB4	15489	34564	0.32%	0.058%
4	4.630	107	110	119	rVB3	14298	32158	0.29%	0.054%
5	4.835	123	128	137	rVV2	24433	63829	0.58%	0.107%
6	5.886	215	220	229	rBV	710338	1345943	12.31%	2.263%
7	7.438	351	356	373	rBV	53587	139778	1.28%	0.235%
8	8.237	421	426	435	rVB	28744	65014	0.59%	0.109%
9	9.710	551	555	561	rVV	3426369	6371379	58.29%	10.714%
10	9.859	561	568	577	rVB	10469	30559	0.28%	0.051%
11	13.192	855	860	865	rBV	4907340	9049049	82.79%	15.216%
12	18.341	1305	1311	1315	rBV	4999343	10039064	91.85%	16.881%
13	22.280	1651	1656	1661	rBV	39670	74636	0.68%	0.126%
14	22.634	1681	1687	1693	rBV2	4908526	10930435	100.00%	18.380%
15	22.771	1693	1699	1709	rVB	53542	174351	1.60%	0.293%
16	27.578	2115	2120	2127	rVB2	15077	37002	0.34%	0.062%
17	27.726	2127	2133	2139	rBV	938827	1658982	15.18%	2.790%
18	30.489	2369	2375	2381	rBV2	4406984	10487239	95.95%	17.635%
19	33.400	2627	2630	2639	rBV2	17663	43957	0.40%	0.074%
20	34.416	2713	2719	2749	rBV	3376706	8813402	80.63%	14.820%

Sum of corrected areas: 59469942

LSC Report - Integrated Chromatogram

File : C:\MSDCHEM\1\5973N\02JAN16\00699.D
 Operator :
 Acquired : 16 Jan 2002 1:20 pm using AcqMethod SCAN508
 Instrument : Instrumen
 Sample Name: 508scan
 Misc Info : January 16, 2002
 Vial Number: 4
 Quant File :SCAN508.RES (RTE Integrator)



Library Search Compound Report

Data File : C:\MSDCHEM\1\5973N\02JAN16\00699.D

Acq On : 16 Jan 2002 1:20 pm

Sample : 508scan

Misc : January 16, 2002

MS Integration Params: LSCINT.P

Vial: 4

Operator:

Inst : Instrumen

Multiplr: 1.00

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

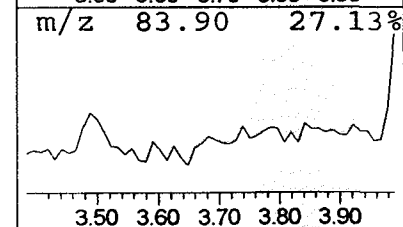
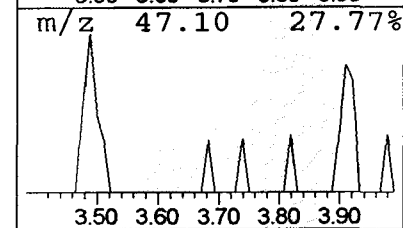
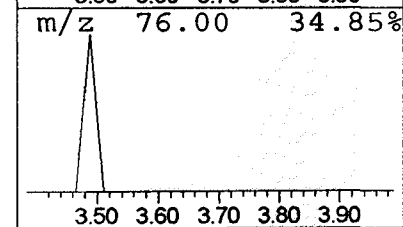
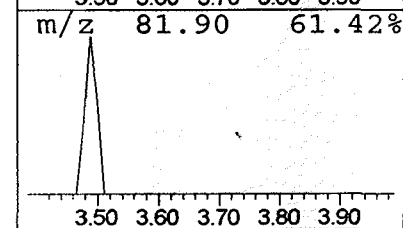
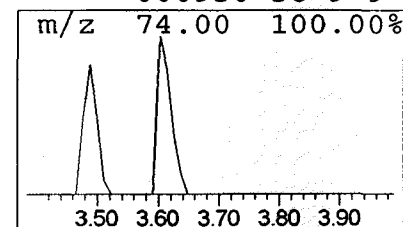
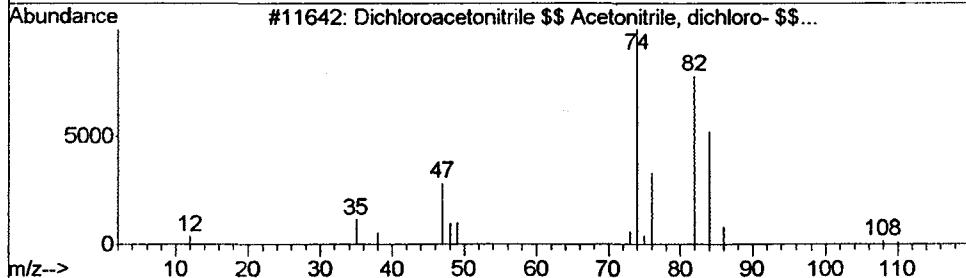
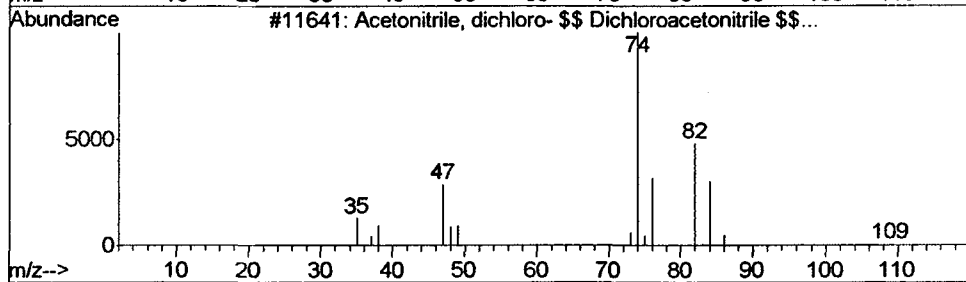
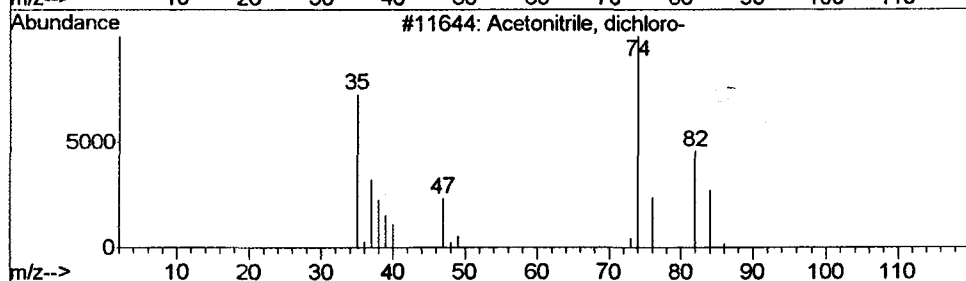
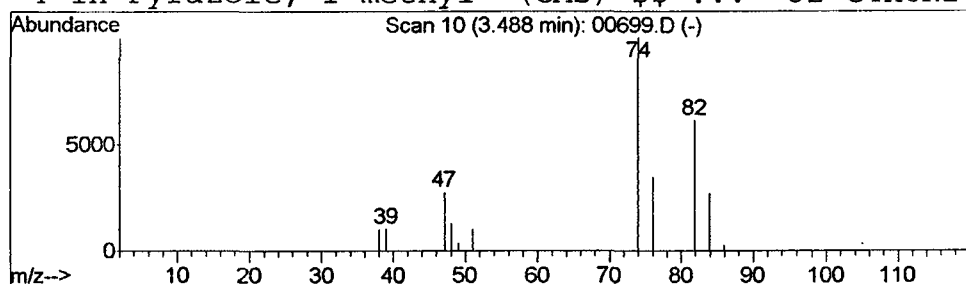
Title : 508 scan method

Library : C:\DATABASE\WILEY7N.L

Peak Number 1 Acetonitrile, dichloro- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.49	0.11 ug/L	33887	1,4-Dichlorobenzene-d4	9.71

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Acetonitrile, dichloro-	109	C2HCl2N	003018-12-0	78
2			Acetonitrile, dichloro- \$\$ Dichl...	109	C2HCl2N	003018-12-0	74
3			Dichloroacetonitrile \$\$ Acetonit...	109	C2HCl2N	003018-12-0	56
4			1H-Pyrazole, 1-methyl- (CAS) \$\$...	82	C4H6N2	000930-36-9	9



Library Search Compound Report

Data File : C:\MSDCHEM\1\5973N\02JAN16\00699.D
 Acq On : 16 Jan 2002 1:20 pm
 Sample : 508scan
 Misc : January 16, 2002
 MS Integration Params: LSCINT.P

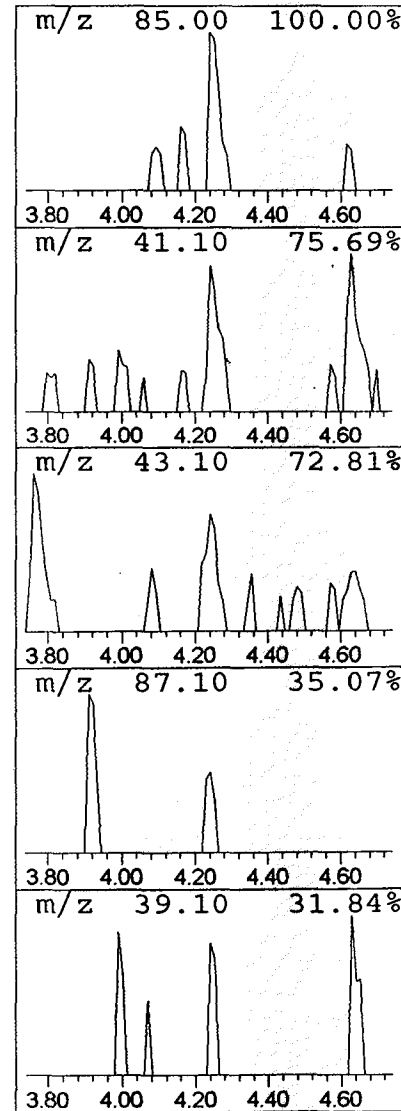
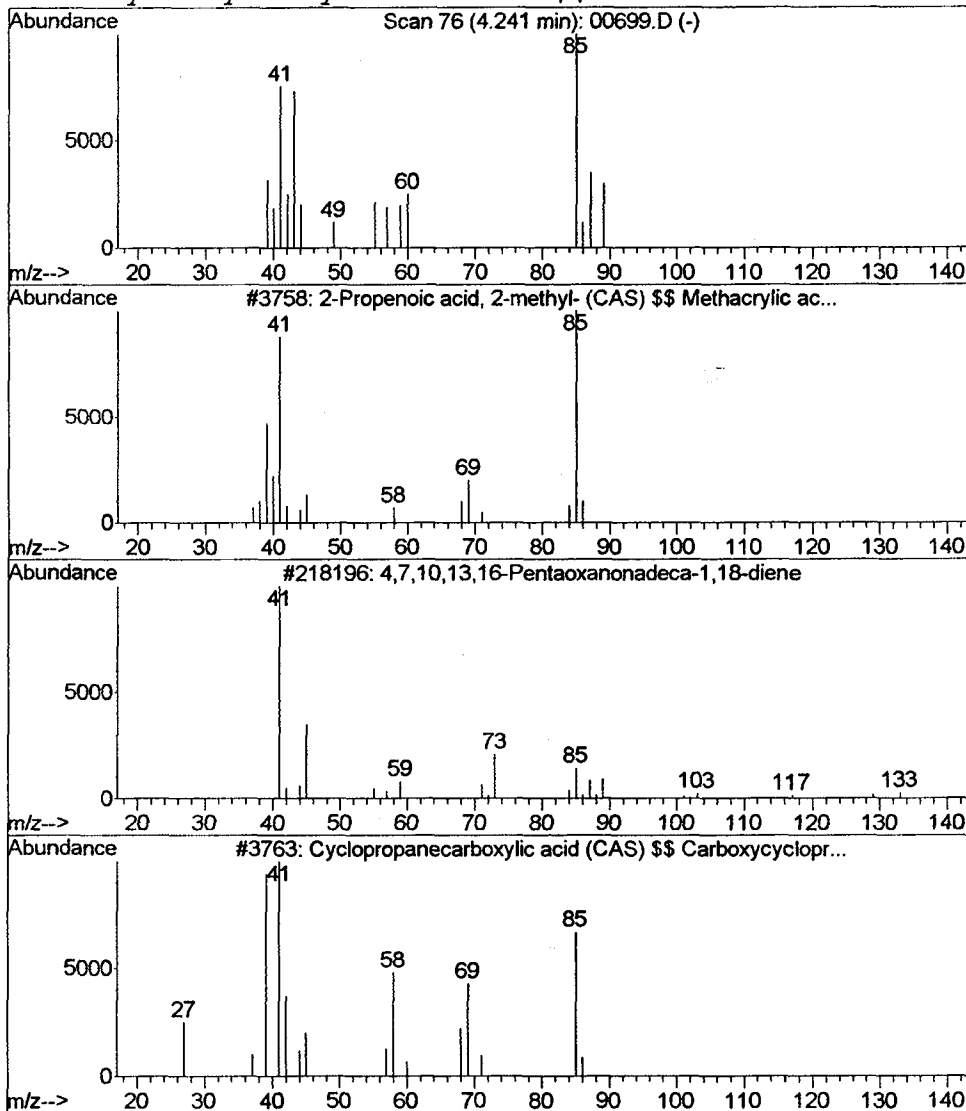
Vial: 4
 Operator:
 Inst : Instrumen
 Multiplr: 1.00

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

 Peak Number 2 2-Propenoic acid, 2-methyl-... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.24	0.11 ug/L	34564	1,4-Dichlorobenzene-d4	9.71

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Propenoic acid, 2-methyl- (CAS...	86	C4H6O2	000079-41-4	30
2			4,7,10,13,16-Pentaoxanonadeca-1,...	274	C14H26O5	058185-54-9	28
3			Cyclopropanecarboxylic acid (CAS...	86	C4H6O2	001759-53-1	25
4			N-Hydroxymethylacetamide \$\$ Acet...	89	C3H7NO2	000625-51-4	10



Library Search Compound Report

Data File : C:\MSDCHEM\1\5973N\02JAN16\00699.D
 Acq On : 16 Jan 2002 1:20 pm
 Sample : 508scan
 Misc : January 16, 2002
 MS Integration Params: LSCINT.P

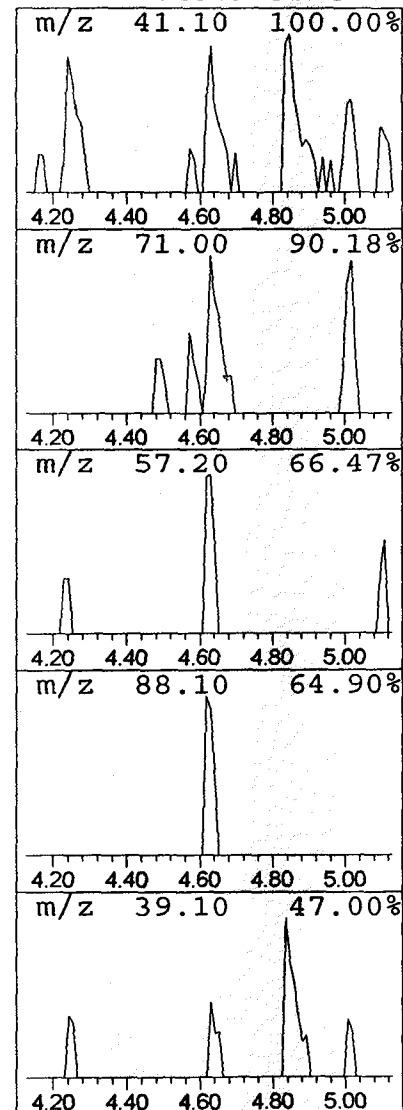
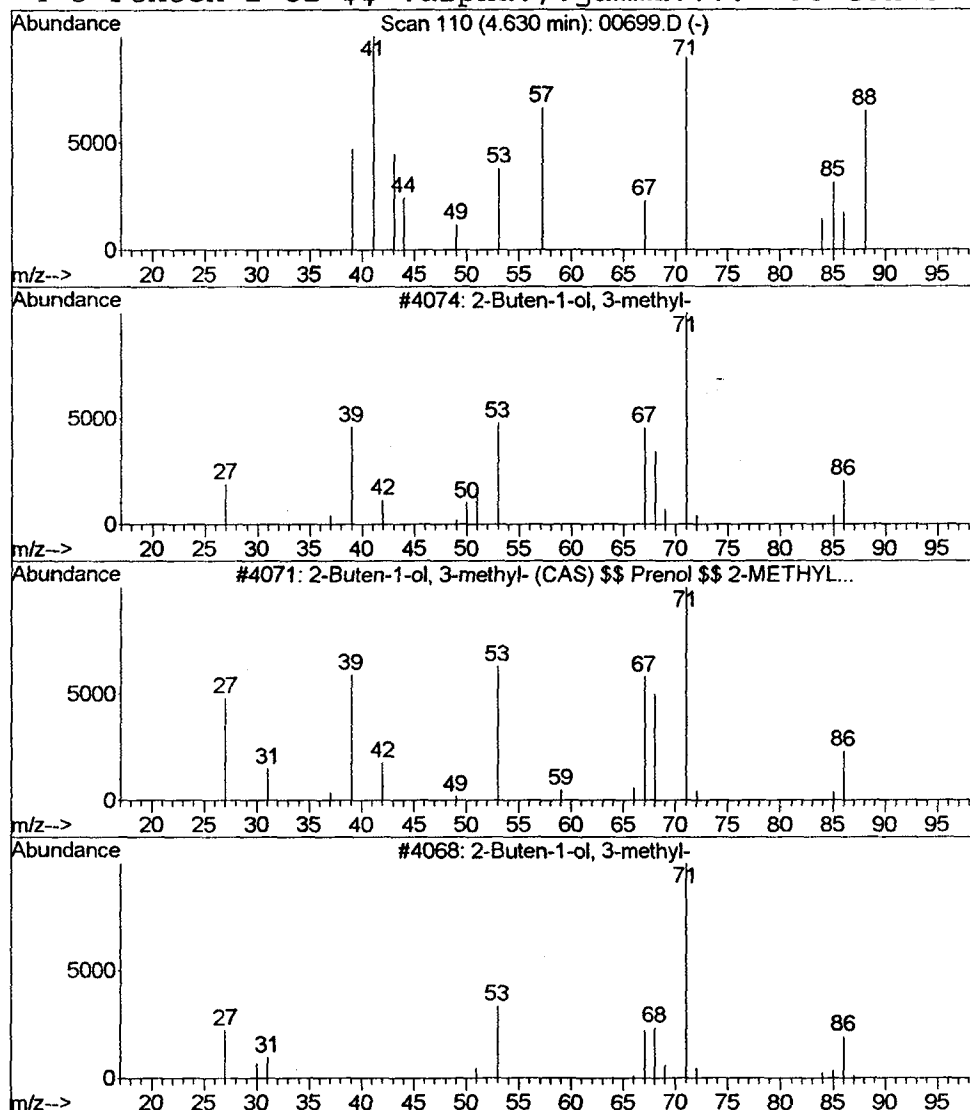
Vial: 4
 Operator:
 Inst : Instrumen
 Multiplr: 1.00

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

 Peak Number 3 2-Buten-1-ol, 3-methyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.63	0.10 ug/L	32158	1,4-Dichlorobenzene-d4	9.71

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Buten-1-ol, 3-methyl-	86	C5H10O	000556-82-1	53
2			2-Buten-1-ol, 3-methyl- (CAS) \$\$...	86	C5H10O	000556-82-1	50
3			2-Buten-1-ol, 3-methyl-	86	C5H10O	000556-82-1	45
4			3-Penten-2-ol \$\$.alpha.,.gamma....	86	C5H10O	001569-50-2	42



Library Search Compound Report

Data File : C:\MSDCHEM\1\5973N\02JAN16\00699.D
 Acq On : 16 Jan 2002 1:20 pm
 Sample : 508scan
 Misc : January 16, 2002
 MS Integration Params: LSCINT.P

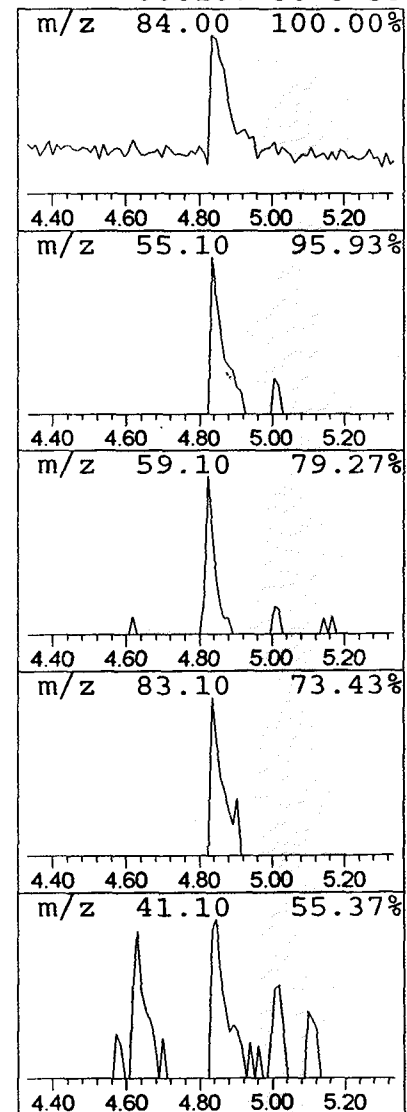
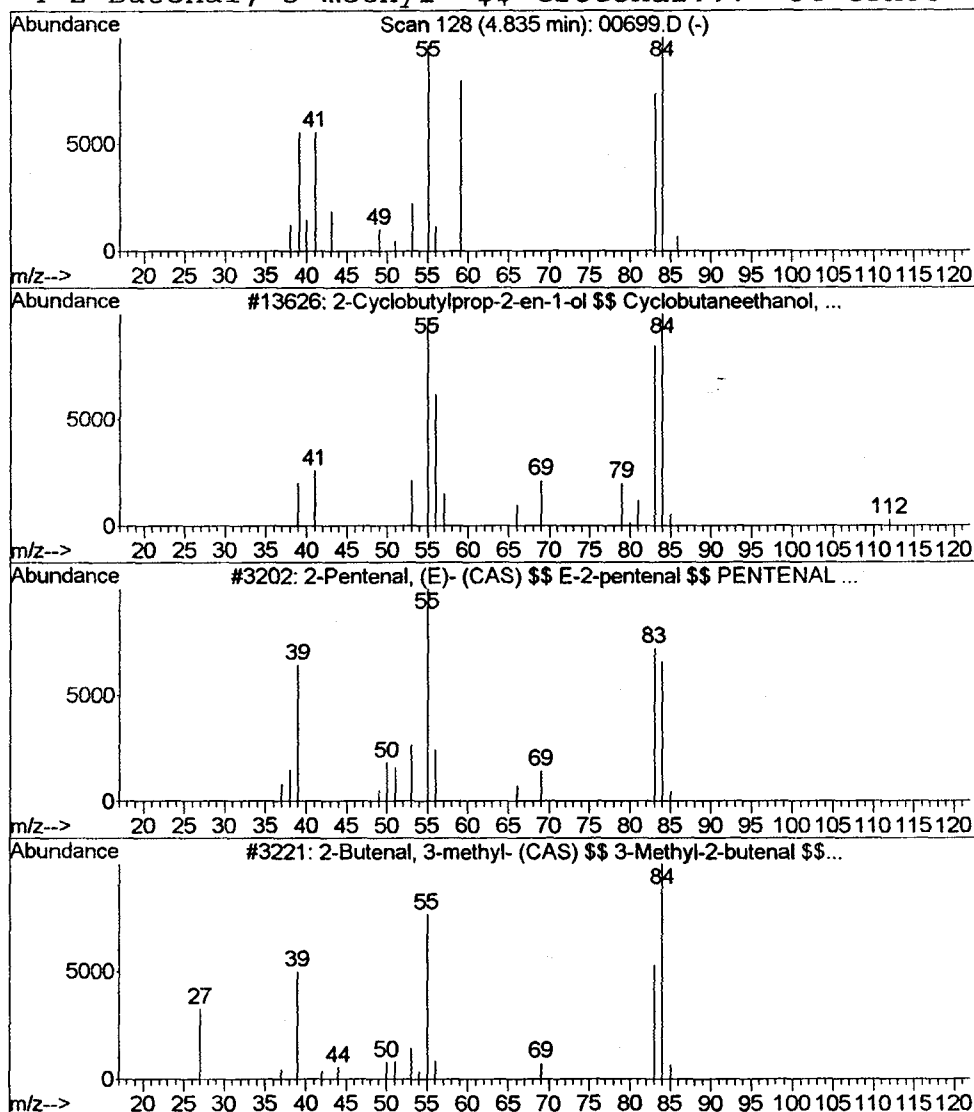
Vial: 4
 Operator:
 Inst : Instrumen
 Multiplr: 1.00

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

 Peak Number 4 2-Cyclobutylprop-2-en-1-ol ... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.84	0.20 ug/L	63829	1,4-Dichlorobenzene-d4	9.71

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Cyclobutylprop-2-en-1-ol \$\$ Cy...	112	C7H12O	116203-80-6	53
2			2-Pentenal, (E)- (CAS) \$\$ E-2-pe...	84	C5H8O	001576-87-0	53
3			2-Butenal, 3-methyl- (CAS) \$\$ 3-...	84	C5H8O	000107-86-8	53
4			2-Butenal, 3-methyl- \$\$ Crotonal...	84	C5H8O	000107-86-8	53



Library Search Compound Report

Data File : C:\MSDCHEM\1\5973N\02JAN16\00699.D
 Acq On : 16 Jan 2002 1:20 pm
 Sample : 508scan
 Misc : January 16, 2002
 MS Integration Params: LSCINT.P

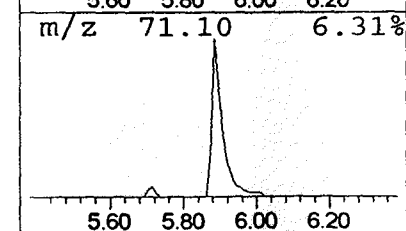
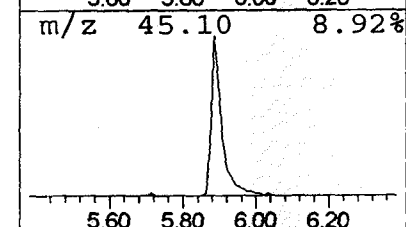
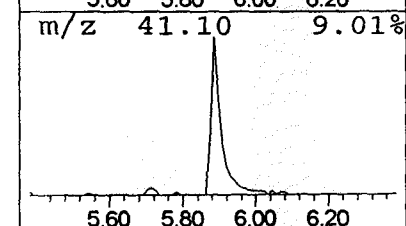
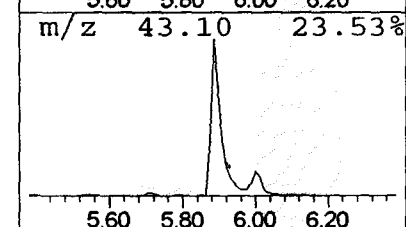
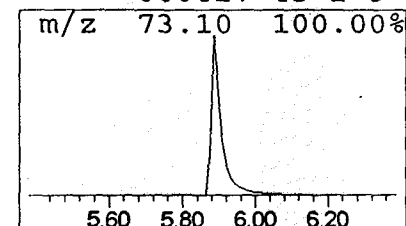
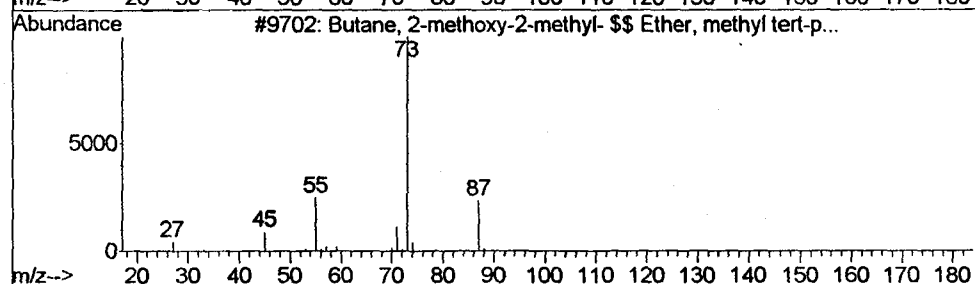
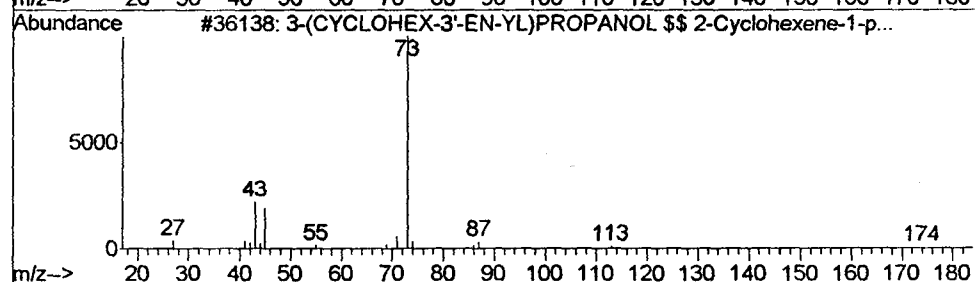
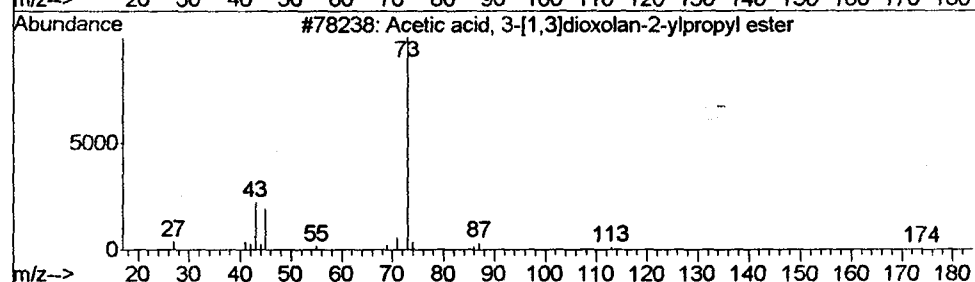
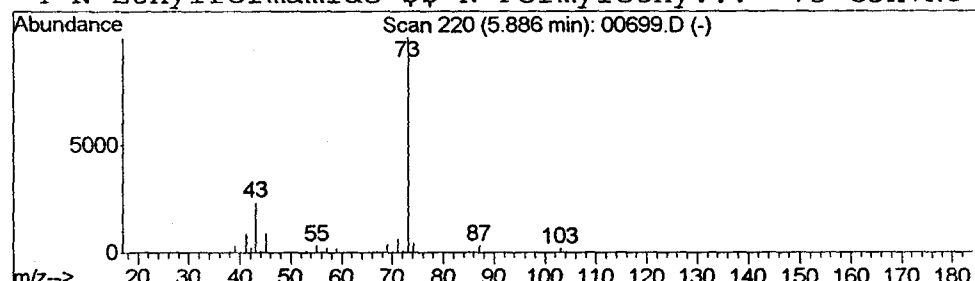
Vial: 4
 Operator:
 Inst : Instrumen
 Multiplr: 1.00

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

 Peak Number 5 Acetic acid, 3-[1,3]dioxola... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.89	4.22 ug/L	1345940	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	36
2		3-(CYCLOHEX-3'-EN-YL)PROPANOL \$\$...	174	C8H14O4	015745-87-6	36
3		Butane, 2-methoxy-2-methyl- \$\$ E...	102	C6H14O	000994-05-8	33
4		N-Ethylformamide \$\$ N-Formylethy...	73	C3H7NO	000627-45-2	9



Library Search Compound Report

Data File : C:\MSDCHEM\1\5973N\02JAN16\00699.D
 Acq On : 16 Jan 2002 1:20 pm
 Sample : 508scan
 Misc : January 16, 2002
 MS Integration Params: LSCINT.P

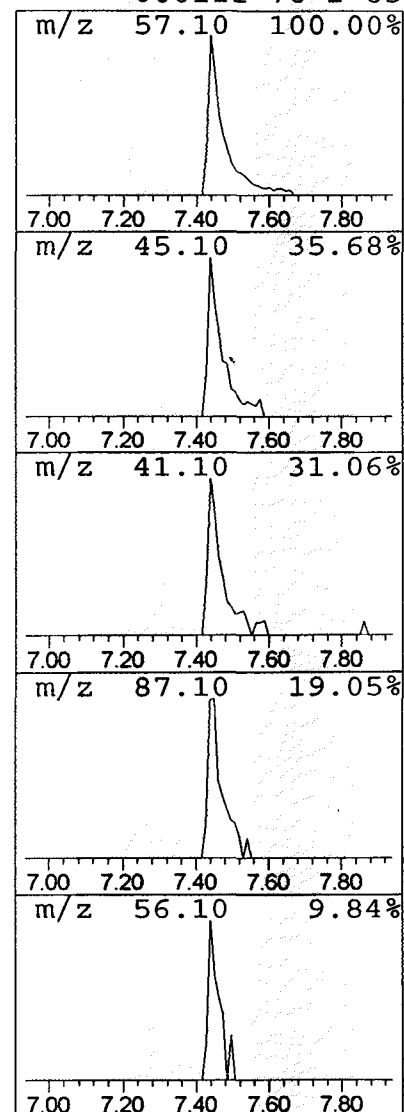
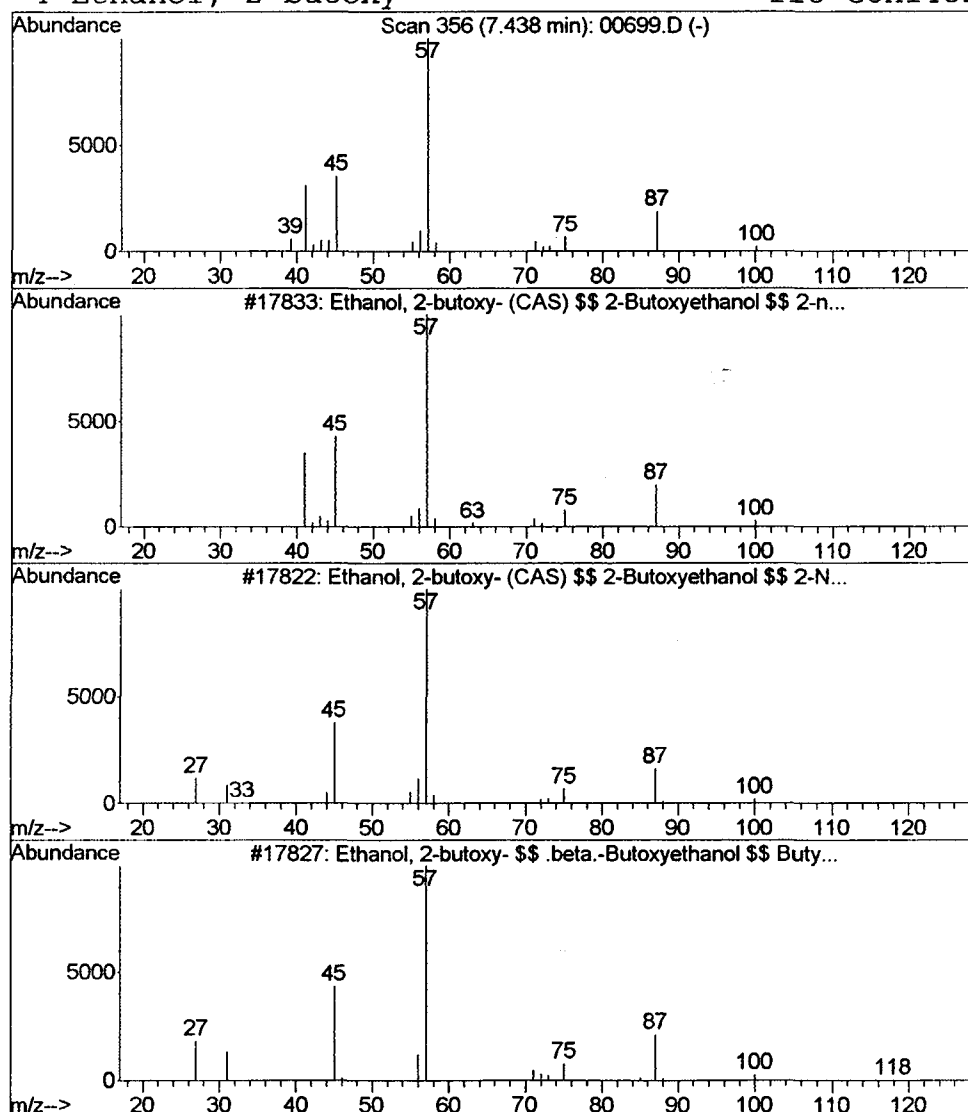
Vial: 4
 Operator:
 Inst : Instrumen
 Multiplr: 1.00

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

 Peak Number 6 Ethanol, 2-butoxy- (CAS) \$\$... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.44	0.44 ug/L	139778	1,4-Dichlorobenzene-d4	9.71

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Ethanol, 2-butoxy- (CAS) \$\$ 2-Bu...	118	C6H14O2	000111-76-2	90
2			Ethanol, 2-butoxy- (CAS) \$\$ 2-Bu...	118	C6H14O2	000111-76-2	83
3			Ethanol, 2-butoxy- \$\$.beta.-But...	118	C6H14O2	000111-76-2	83
4			Ethanol, 2-butoxy-	118	C6H14O2	000111-76-2	83



Library Search Compound Report

Data File : C:\MSDCHEM\1\5973N\02JAN16\00699.D
Acq On : 16 Jan 2002 1:20 pm
Sample : 508scan
Misc : January 16, 2002
MS Integration Params: LSCINT.P

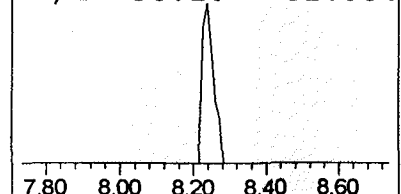
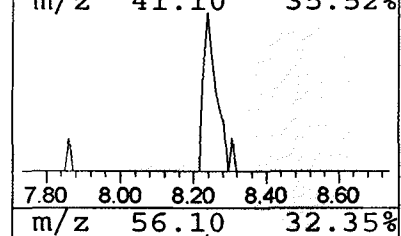
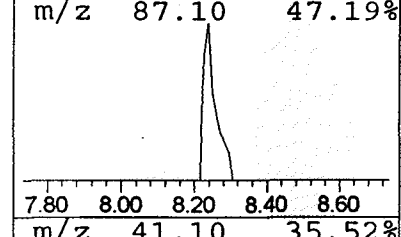
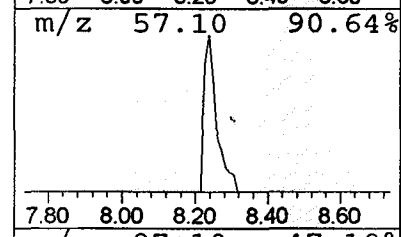
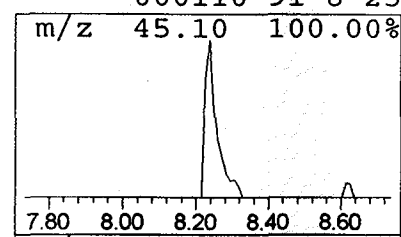
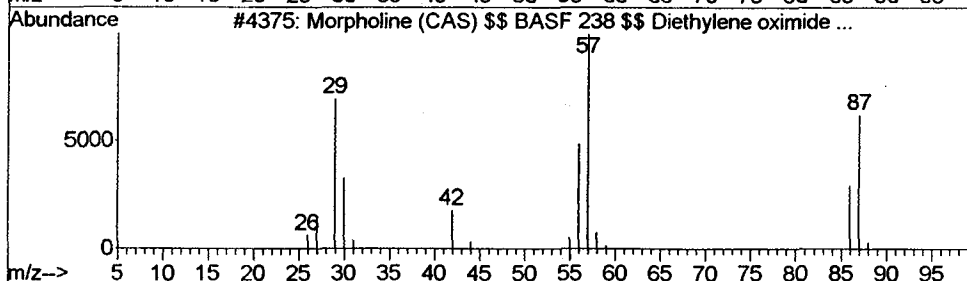
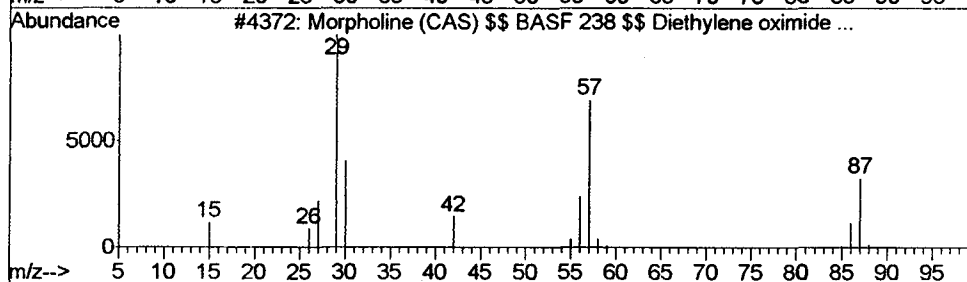
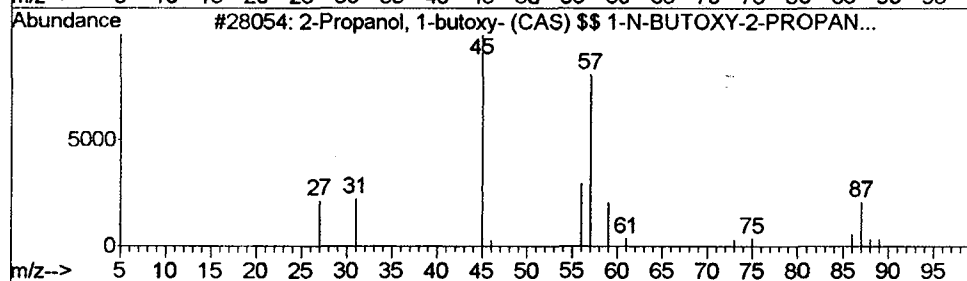
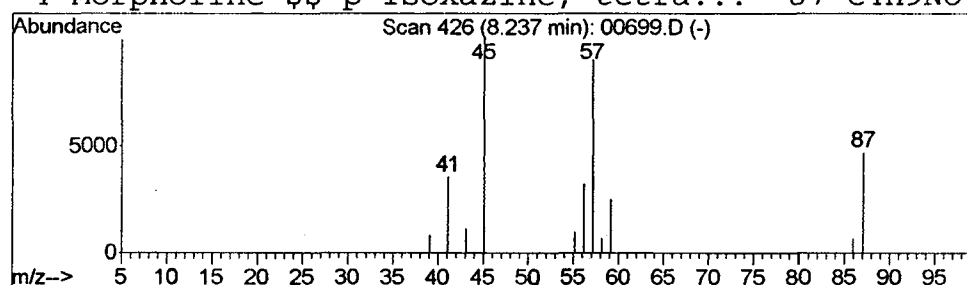
Vial: 4
Operator:
Inst : Instrumen
Multiplr: 1.00

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

Peak Number 7 2-Propanol, 1-butoxy- (CAS)... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.24	0.20 ug/L	65014	1,4-Dichlorobenzene-d4	9.71

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Propanol, 1-butoxy- (CAS) \$\$ 1...	132	C7H16O2	005131-66-8	56
2		Morpholine (CAS) \$\$ BASF 238 \$\$...	87	C4H9NO	000110-91-8	35
3		Morpholine (CAS) \$\$ BASF 238 \$\$...	87	C4H9NO	000110-91-8	25
4		Morpholine \$\$ p-Isoxazine, tetra...	87	C4H9NO	000110-91-8	25



Library Search Compound Report

Data File : C:\MSDCHEM\1\5973N\02JAN16\00699.D

Acq On : 16 Jan 2002 1:20 pm

Sample : 508scan

Misc : January 16, 2002

MS Integration Params: LSCINT.P

Vial: 4

Operator:

Inst : Instrumen

Multiplr: 1.00

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

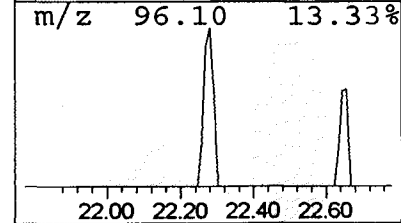
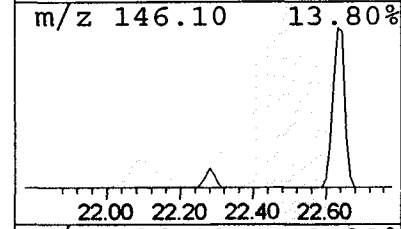
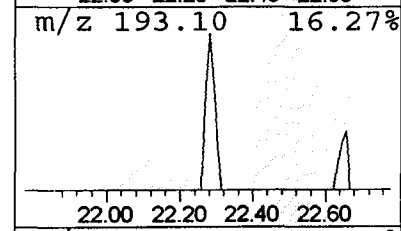
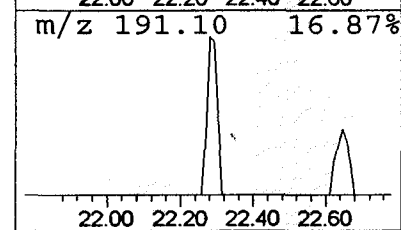
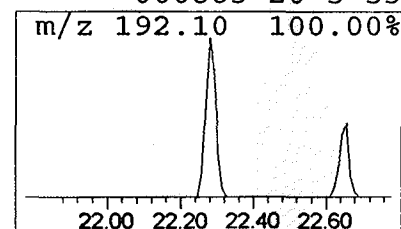
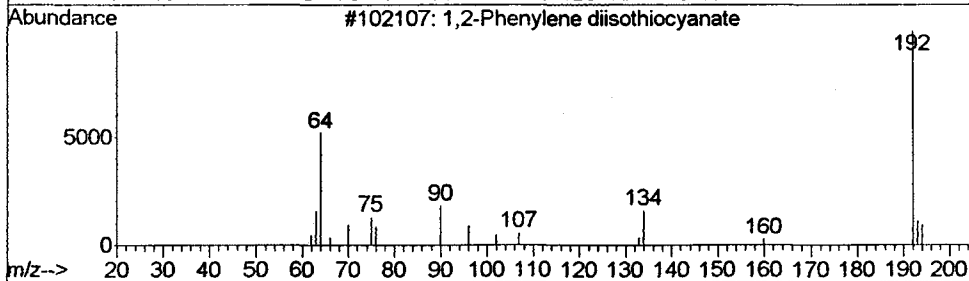
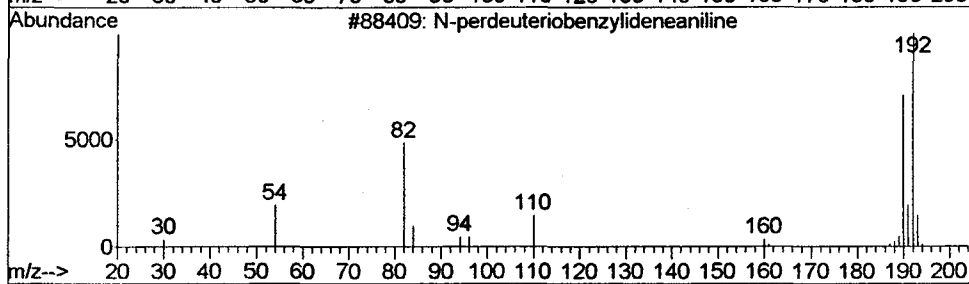
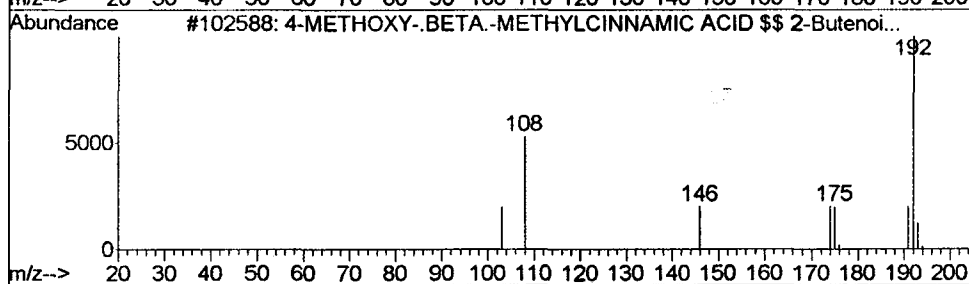
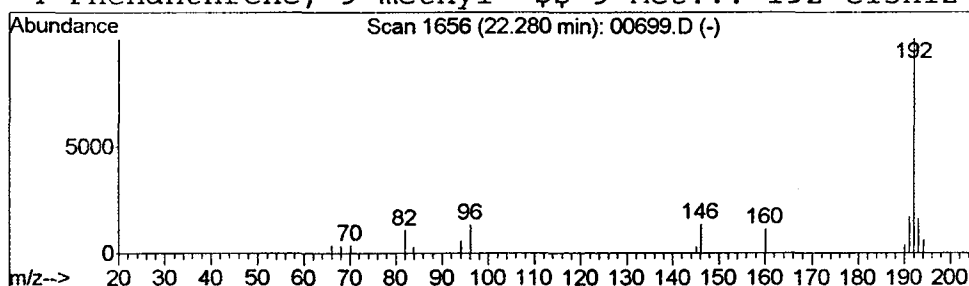
Title : 508 scan method

Library : C:\DATABASE\WILEY7N.L

Peak Number 8 4-METHOXY-.BETA.-METHYLCINN... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.28	0.14 ug/L	74636	Phenanthrene-d10	22.63

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		4-METHOXY-.BETA.-METHYLCINNAMIC ...	192	C11H12O3	019169-94-9	72
2		N-perdeuteriobenzylideneaniline	192	C13D11N	000538-51-2	59
3		1,2-Phenylene diisothiocyanate	192	C8H4N2S2	071105-17-4	53
4		Phenanthrene, 9-methyl- \$\$ 9-Met...	192	C15H12	000883-20-5	53



Library Search Compound Report

Data File : C:\MSDCHEM\1\5973N\02JAN16\00699.D
 Acq On : 16 Jan 2002 1:20 pm
 Sample : 508scan
 Misc : January 16, 2002
 MS Integration Params: LSCINT.P

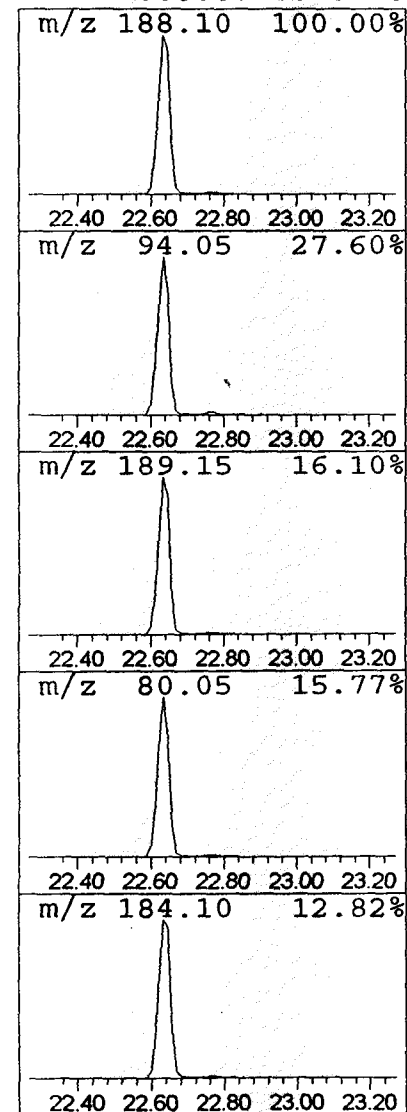
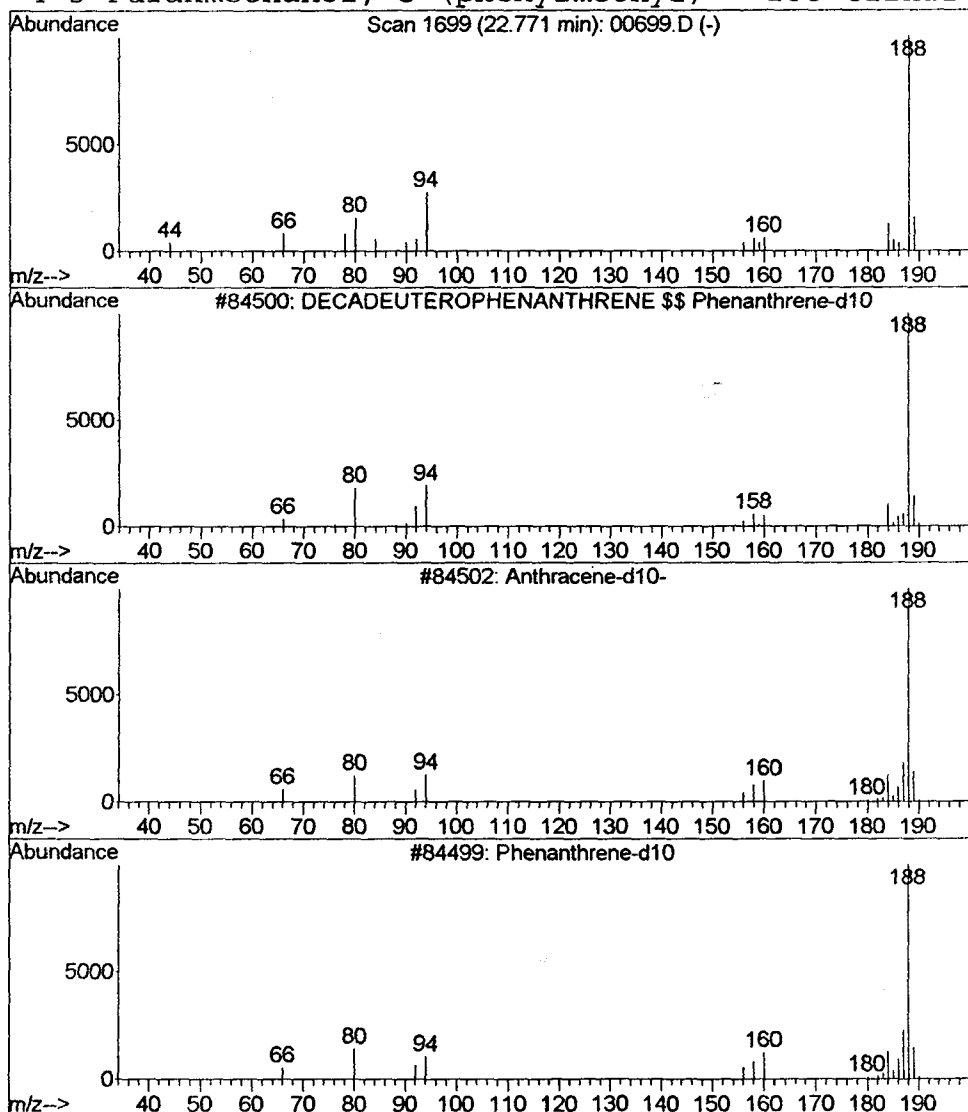
Vial: 4
 Operator:
 Inst : Instrumen
 Multiplr: 1.00

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

 Peak Number 9 DECADEUTEROPHENANTHRENE \$\$... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.77	0.32 ug/L	174351	Phenanthrene-d10	22.63

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	DECADEUTEROPHENANTHRENE \$\$ Phena...	188	C14D10	001517-22-2	78
2		Anthracene-d10-	188	C14D10	001719-06-8	64
3		Phenanthrene-d10	188	C14D10	001517-22-2	64
4		3-Furanmethanol, 5-(phenylmethyl)-	188	C12H12O2	000000-00-0	50



Tentatively Identified Compound (LSC) summary

Operator ID: Date Acquired: 16 Jan 2002 1:20 pm
 Data File: C:\MSDCHEM\1\5973N\02JAN16\00699.D
 Name: 508scan
 Date: January 16, 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...	3.49	0.1	ug/L	33887	1	9.71	6371380	20.0
2-Propenoic acid,...	4.24	0.1	ug/L	34564	1	9.71	6371380	20.0
2-Buten-1-ol, 3-m...	4.63	0.1	ug/L	32158	1	9.71	6371380	20.0
2-Cyclobutylprop...	4.84	0.2	ug/L	63829	1	9.71	6371380	20.0
Acetic acid, 3-[1...	5.89	4.2	ug/L	1345940	1	9.71	6371380	20.0
Ethanol, 2-butoxy...	7.44	0.4	ug/L	139778	1	9.71	6371380	20.0
2-Propanol, 1-but...	8.24	0.2	ug/L	65014	1	9.71	6371380	20.0
4-METHOXY-.BETA.-...	22.28	0.1	ug/L	74636	4	22.63	10930400	20.0
DECADEUTEROPHENAN...	22.77	0.3	ug/L	174351	4	22.63	10930400	20.0