

User: Vinci, David L  
 Westchester Dept. of Labs & Research  
 Date: 03/05/2002 Time: 12:41:06

Sample: AE02720  
 Analysis: \$GCMS GCMS Scan For Unknowns  
 Units: ug  
 Started: 3/5/02 12:23 Ended: 3/5/02 12:23 Analyst: DLV  
 Result source: manual entry  
 Page: 1

|   | Component Name          | Viol | Result | Qualifier | MDL |
|---|-------------------------|------|--------|-----------|-----|
| 1 | Other unknown compounds |      | .      |           |     |
| 2 | Surr_2,4-DDD            |      | 113 ✓  |           | 5.0 |

2/4/2  
 Amciv

Sent 3/5/2  
 with Data-MS

**Comments for Sample AE02720 - Analysis \$GCMS**

Westchester Dept. of Labs & Research

User: Vinci, David L

Date: 03-05-2002 Time: 12:41:02

The GCMS scan, from 35 to 550 amu, reveals the presence of several compounds, such as 2,2-Dimethoxybutane at 0.44 ug/L with a fit of 83 out of 100; 9H-Carbazole at 0.18 ug/L with a fit of 90 out of 100; n-Heptadecanol at 0.75 ug/L with a fit of 94 out of 100; and Pyrene at 0.14 ug/L and with a fit of 90 out of 100. The recoveries for 14 of the 43 compounds in the LCS Standard were above their acceptance ranges.

## Quantitation Report (Not Reviewed)

Data File : C:\MSDCHEM\1\5973N\02FEB28\02720.D Vial: 5  
Acq On : 28 Feb 2002 6:00 pm Operator:  
Sample : 508 scan Inst : Instrumen  
Misc : February 28, 2002 Multiplr: 1.00  
MS Integration Params: SPLIT.P  
Quant Time: Mar 01 09:03:00 2002 Quant Results File: HP022102.RES

Quant Method : C:\MSDCHEM\1...\HP022102.M (RTE Integrator)  
Title : 508 scan method  
Last Update : Mon Feb 25 10:47:59 2002  
Response via : Initial Calibration  
DataAcq Meth : HP020702

| Internal Standards        | R.T.  | QIon | Response | Conc  | Units | Dev (Min) |
|---------------------------|-------|------|----------|-------|-------|-----------|
| 1) 1,4-Dichlorobenzene-d4 | 9.71  | 152  | 1518931  | 20.00 | ug/L  | 0.00      |
| 10) Naphthalene-d8        | 13.20 | 136  | 6594257  | 20.00 | ug/L  | 0.00      |
| 17) Acenaphthene-d10      | 18.34 | 164  | 3317757  | 20.00 | ug/L  | 0.00      |
| 29) Phenanthrene-d10      | 22.63 | 188  | 5556003  | 20.00 | ug/L  | -0.01     |
| 38) Chrysene-d12          | 30.49 | 240  | 4631181  | 20.00 | ug/L  | -0.03     |
| 44) Perylene-d12          | 34.42 | 264  | 3264987  | 20.00 | ug/L  | -0.03     |

System Monitoring Compounds

|               |       |     |          |      |         |      |
|---------------|-------|-----|----------|------|---------|------|
| 37) 2,4-DDD   | 27.72 | 235 | 418136   | 5.65 | ug/L    | 0.00 |
| Spiked Amount | 5.000 |     | Recovery | =    | 113.00% |      |

Target Compounds

|                         |       |     |        |      |      | Qvalue |
|-------------------------|-------|-----|--------|------|------|--------|
| 20) Dimethylphthalate   | 18.34 | 163 | 535753 | 2.66 | ug/L | # 58   |
| 21) 2,6-Dinitrotoluene  | 18.33 | 165 | 427566 | 9.20 | ug/L | # 27   |
| 35) Di-n-butylphthalate | 24.86 | 149 | 146882 | 0.41 | ug/L | 97     |

*Amcis 2/4/2*

Quantitation Report

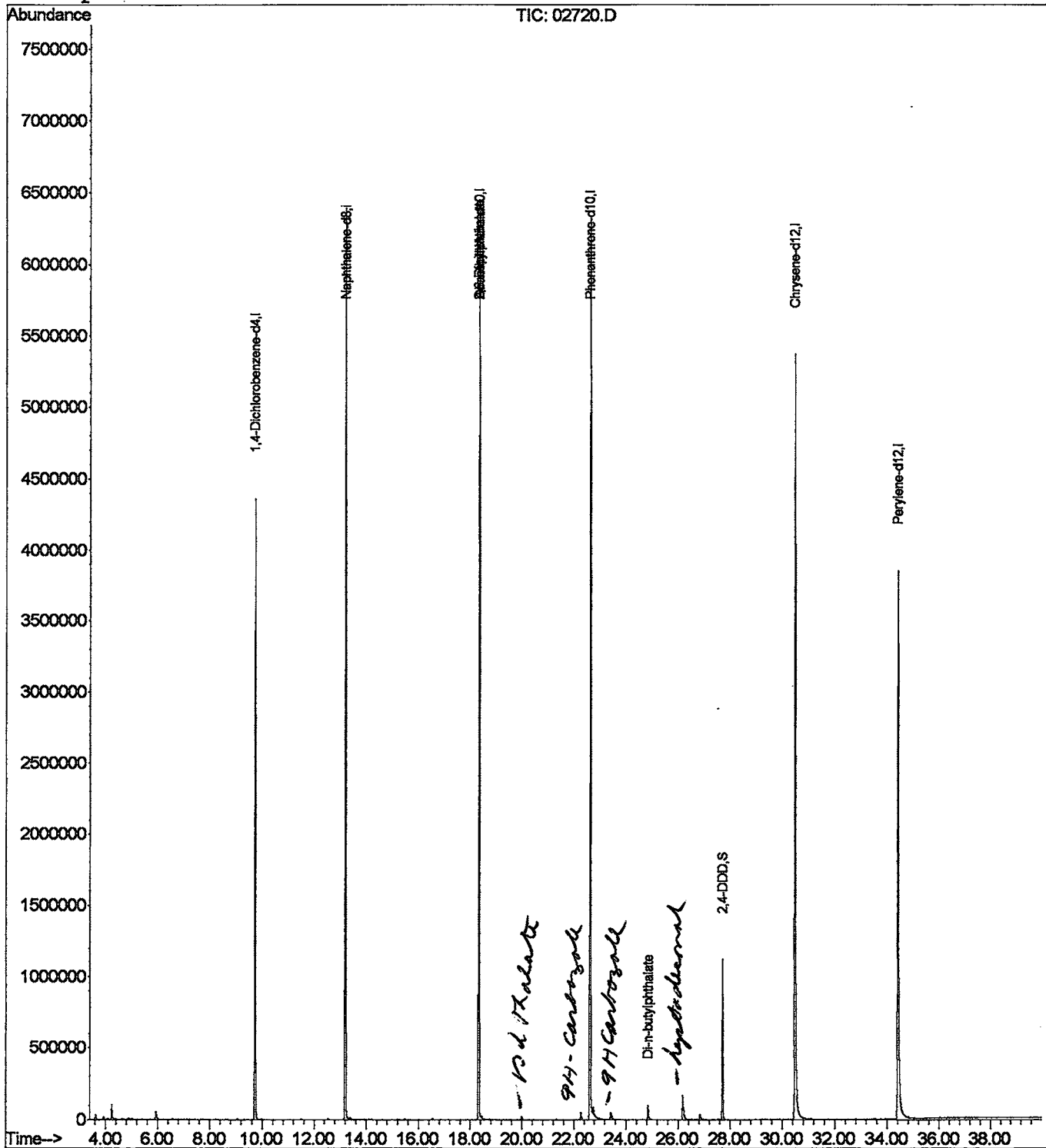
(Not Reviewed)

Data File : C:\MSDCHEM\1\5973N\02FEB28\02720.D  
 Acq On : 28 Feb 2002 6:00 pm  
 Sample : 508 scan  
 Misc : February 28, 2002  
 MS Integration Params: SPLIT.P  
 Quant Time: Mar 1 9:03 2002

Vial: 5  
 Operator:  
 Inst : Instrumen  
 Multiplr: 1.00

Quant Results File: HP022102.R

Method : C:\MSDCHEM\1\METHODS\METHODS\HP022102.M (RTE Integrator)  
 Title : 508 scan method  
 Last Update : Mon Feb 25 10:47:59 2002  
 Response via : Initial Calibration



## LSC Area Percent Report

Data File : C:\MSDCHEM\1\5973N\02FEB28\02720.D  
 Acq On : 28 Feb 2002 6:00 pm  
 Sample : 508 scan  
 Misc : February 28, 2002  
 MS Integration Params: LSCINT.P

Vial: 5  
 Operator:  
 Inst : Instrumen  
 Multiplr: 1.00

Method : C:\MSDCHEM\1\METHODS\METHODS\HP022102.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

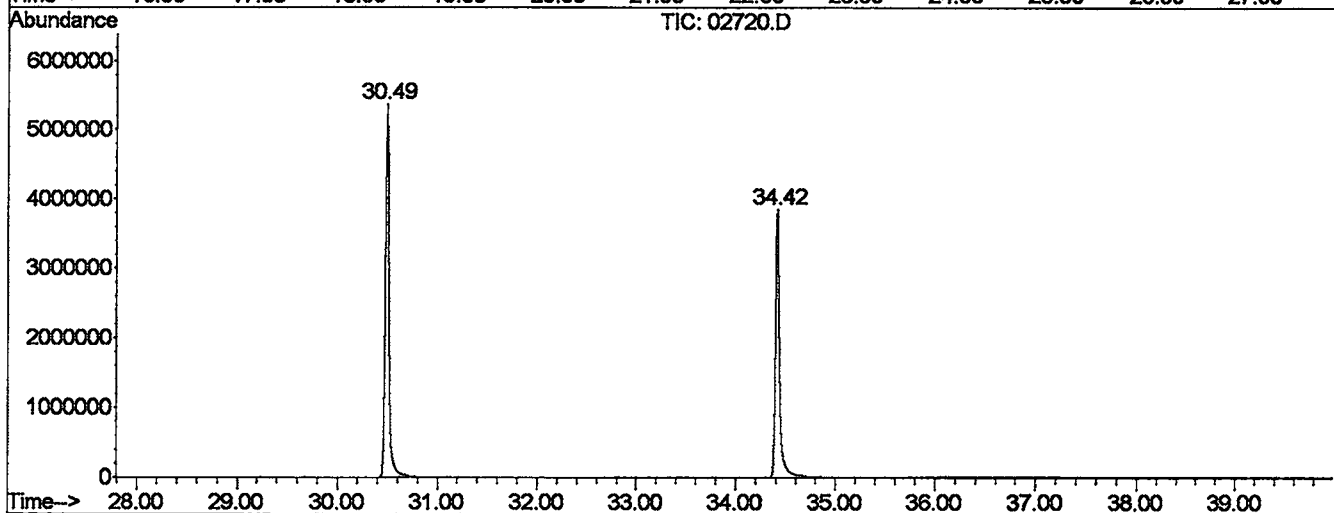
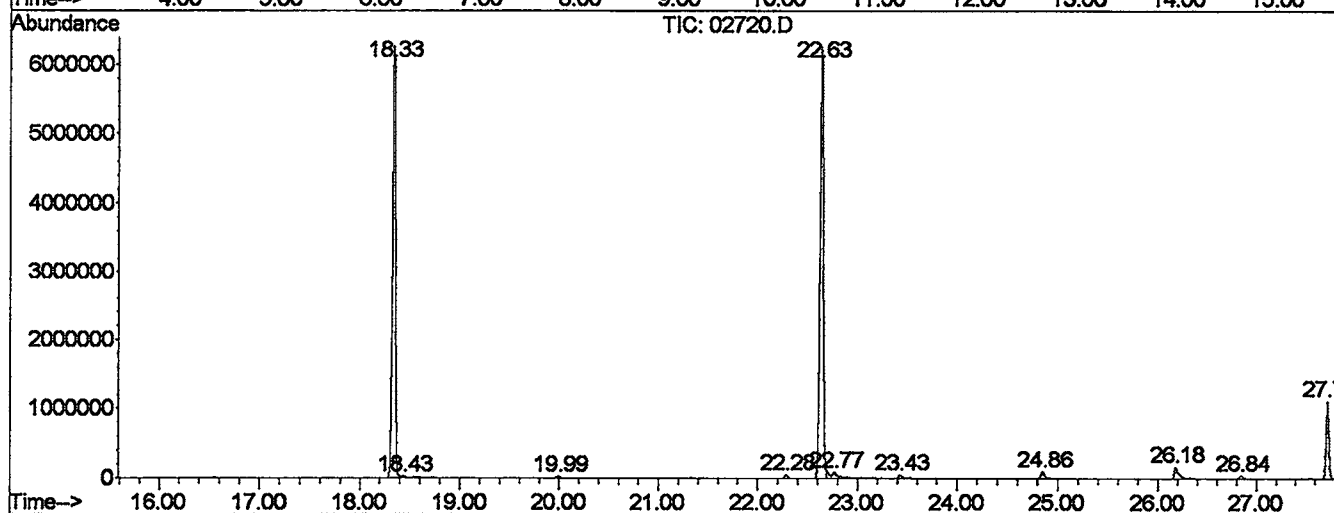
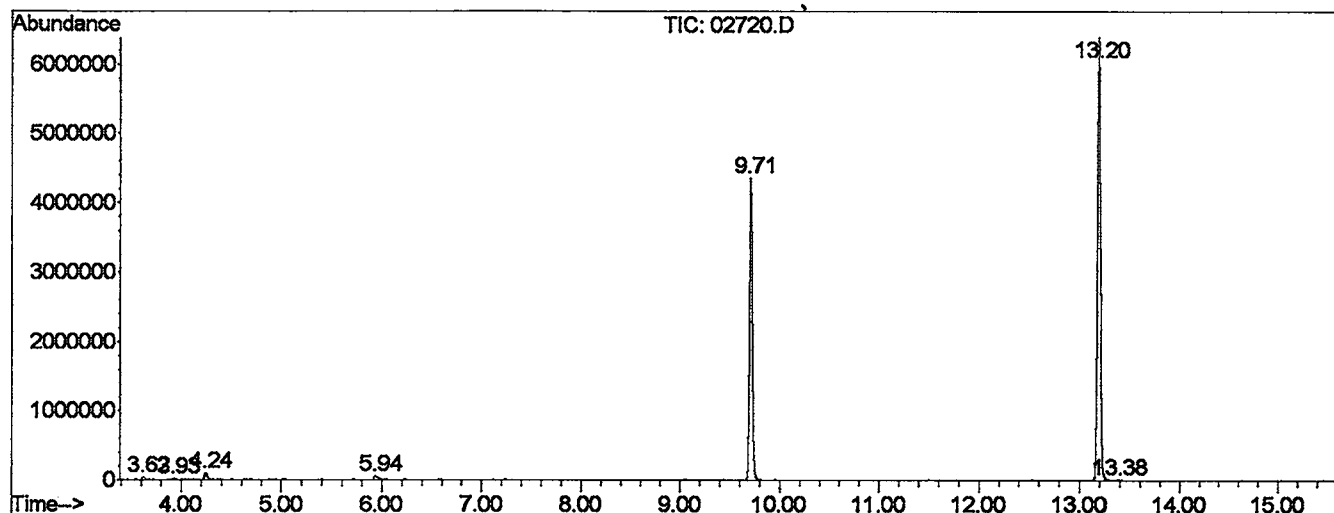
Signal : TIC

| peak<br># | R.T.<br>min | first<br>scan | max<br>scan | last<br>scan | PK<br>TY | peak<br>height | corr.<br>area | corr.<br>% max. | % of<br>total |
|-----------|-------------|---------------|-------------|--------------|----------|----------------|---------------|-----------------|---------------|
| 1         | 3.622       | 19            | 22          | 29           | rVB      | 35378          | 65579         | 0.45%           | 0.084%        |
| 2         | 3.927       | 47            | 49          | 53           | rBV      | 19579          | 29091         | 0.20%           | 0.037%        |
| 3         | 4.243       | 73            | 77          | 85           | rVB2     | 106227         | 188579        | 1.30%           | 0.242%        |
| 4         | 5.936       | 225           | 227         | 239          | rBV      | 58968          | 160391        | 1.11%           | 0.206%        |
| 5         | 9.707       | 557           | 561         | 567          | rBV      | 4360252        | 8513116       | 58.75%          | 10.917%       |
| 6         | 13.195      | 865           | 870         | 875          | rBV      | 6389899        | 12479804      | 86.13%          | 16.004%       |
| 7         | 13.376      | 883           | 886         | 893          | rVB2     | 12661          | 37307         | 0.26%           | 0.048%        |
| 8         | 18.331      | 1319          | 1325        | 1331         | rBV      | 6266598        | 13588298      | 93.78%          | 17.425%       |
| 9         | 18.433      | 1331          | 1334        | 1345         | rVB2     | 22290          | 66680         | 0.46%           | 0.086%        |
| 10        | 19.991      | 1467          | 1472        | 1477         | rVV2     | 23183          | 52230         | 0.36%           | 0.067%        |
| 11        | 22.283      | 1669          | 1675        | 1681         | rBV      | 53797          | 101679        | 0.70%           | 0.130%        |
| 12        | 22.633      | 1701          | 1706        | 1711         | rBV2     | 6250307        | 14490062      | 100.00%         | 18.582%       |
| 13        | 22.768      | 1715          | 1718        | 1749         | rVB2     | 85049          | 387292        | 2.67%           | 0.497%        |
| 14        | 23.434      | 1773          | 1777        | 1785         | rBV      | 47705          | 129397        | 0.89%           | 0.166%        |
| 15        | 24.856      | 1899          | 1903        | 1911         | rBV      | 98513          | 230299        | 1.59%           | 0.295%        |
| 16        | 26.177      | 2015          | 2020        | 2035         | rBV2     | 168840         | 541745        | 3.74%           | 0.695%        |
| 17        | 26.843      | 2075          | 2079        | 2089         | rVV      | 38190          | 96136         | 0.66%           | 0.123%        |
| 18        | 27.724      | 2153          | 2157        | 2163         | rBV      | 1120511        | 2015617       | 13.91%          | 2.585%        |
| 19        | 30.490      | 2397          | 2402        | 2435         | rVB2     | 5364756        | 13760187      | 94.96%          | 17.646%       |
| 20        | 34.418      | 2743          | 2750        | 2783         | rBV      | 3850395        | 11047176      | 76.24%          | 14.167%       |

Sum of corrected areas: 77980665

# LSC Report - Integrated Chromatogram

File : C:\MSDCHEM\1\5973N\02FEB28\02720.D  
 Operator :  
 Acquired : 28 Feb 2002 6:00 pm using AcqMethod HP020702  
 Instrument : Instrumen  
 Sample Name: 508 scan  
 Misc Info : February 28, 2002  
 Vial Number: 5  
 Quant File :HP022102.RES (RTE Integrator)



## Library Search Compound Report

Data File : C:\MSDCHEM\1\5973N\02FEB28\02720.D

Acq On : 28 Feb 2002 6:00 pm

Sample : 508 scan

Misc : February 28, 2002

MS Integration Params: LSCINT.P

Vial: 5

Operator:

Inst : Instrumen

Multiplr: 1.00

Quant Method : C:\MSDCHEM\1...\HP022102.M (RTE Integrator)

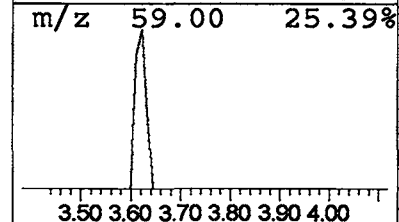
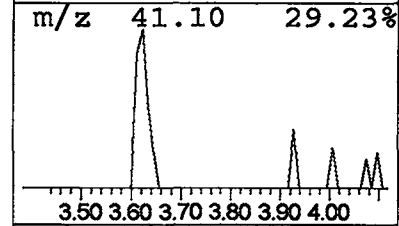
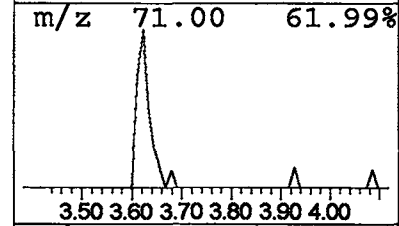
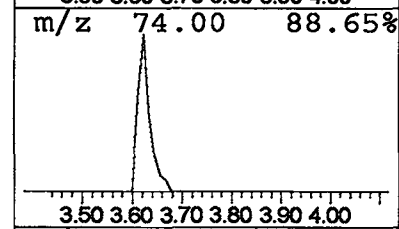
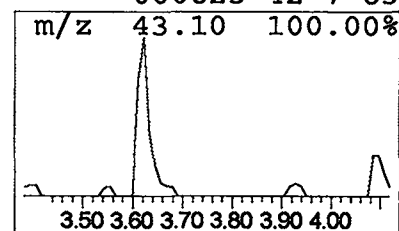
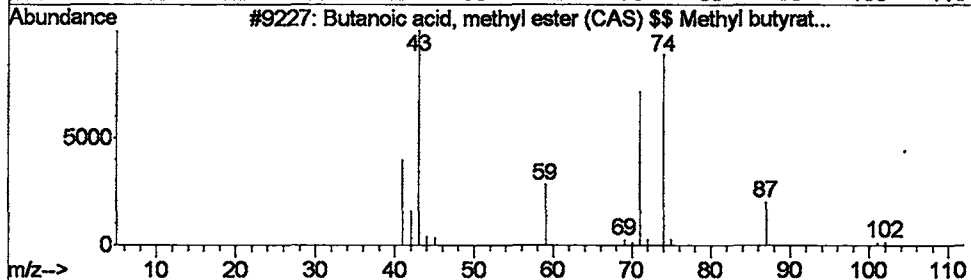
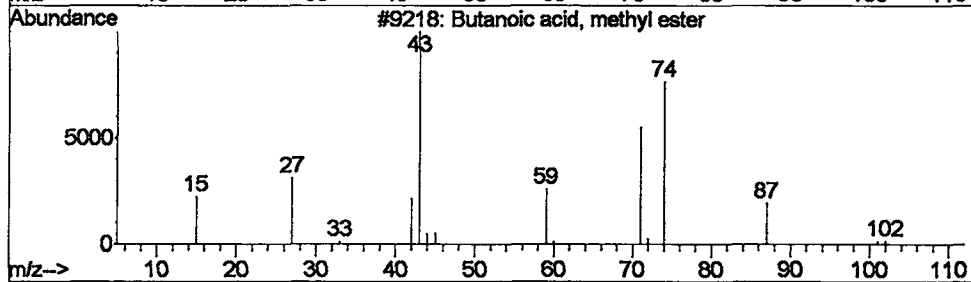
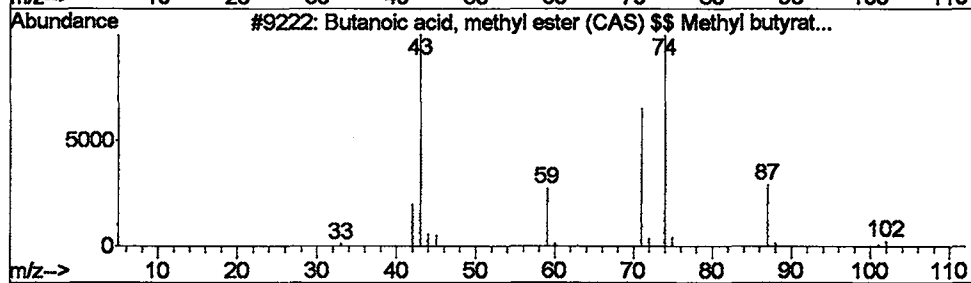
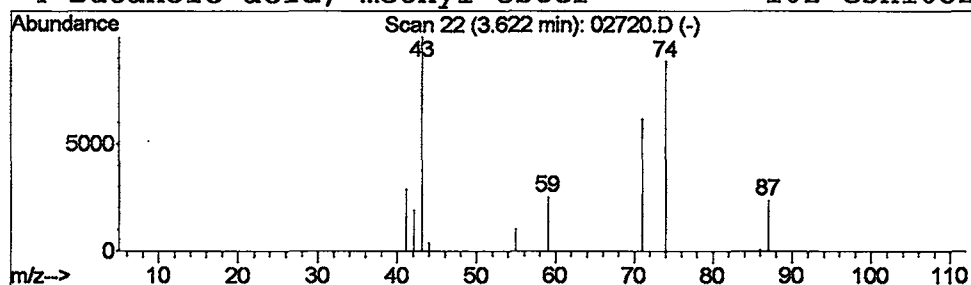
Title : 508 scan method

Library : C:\DATABASE\WILEY7N.L

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

| R.T. | EstConc   | Area  | Relative to ISTD       | R.T. |
|------|-----------|-------|------------------------|------|
| 3.62 | 0.15 ug/L | 65579 | 1,4-Dichlorobenzene-d4 | 9.71 |

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



## Library Search Compound Report

Data File : C:\MSDCHEM\1\5973N\02FEB28\02720.D  
Acq On : 28 Feb 2002 6:00 pm  
Sample : 508 scan  
Misc : February 28, 2002  
MS Integration Params: LSCINT.P

Vial: 5  
Operator:  
Inst : Instrumen  
Multiplr: 1.00

Quant Method : C:\MSDCHEM\1...\HP022102.M (RTE Integrator)

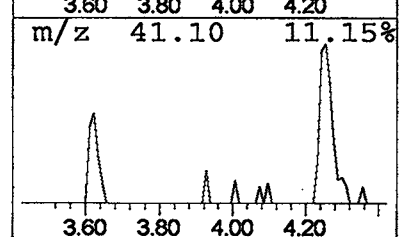
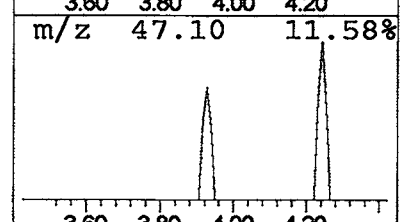
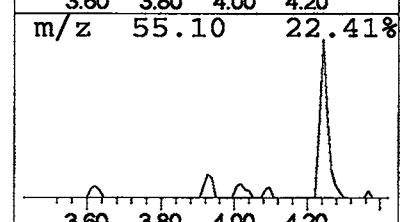
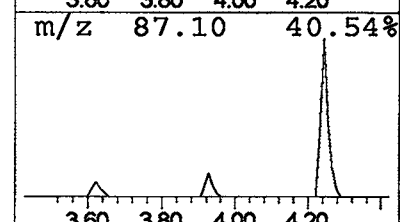
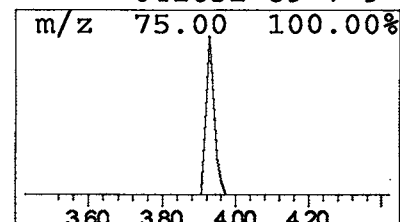
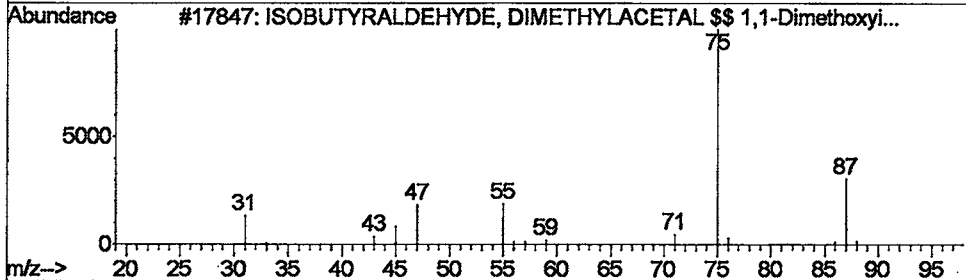
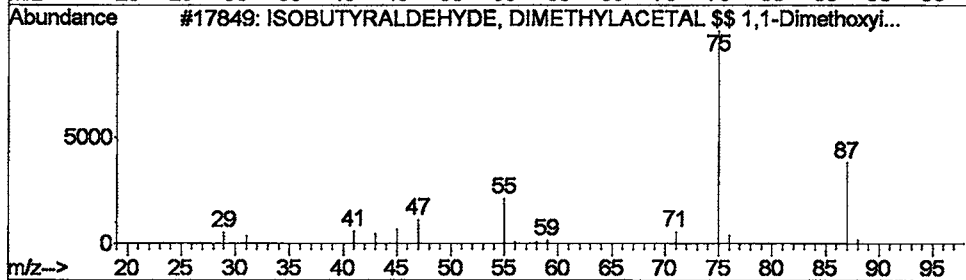
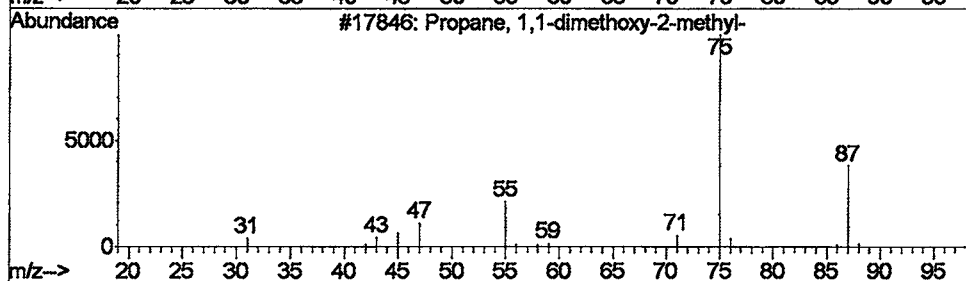
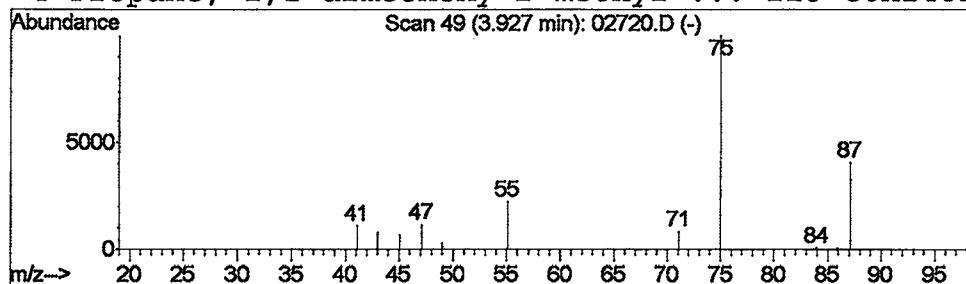
Title : 508 scan method

Library : C:\DATABASE\WILEY7N.L

\*\*\*\*\*  
Peak Number 2 Propane, 1,1-dimethoxy-2-me... Concentration Rank 11

| R.T. | EstConc   | Area  | Relative to ISTD       | R.T. |
|------|-----------|-------|------------------------|------|
| 3.93 | 0.07 ug/L | 29091 | 1,4-Dichlorobenzene-d4 | 9.71 |

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



## Library Search Compound Report

Data File : C:\MSDCHEM\1\5973N\02FEB28\02720.D  
Acq On : 28 Feb 2002 6:00 pm  
Sample : 508 scan  
Misc : February 28, 2002  
MS Integration Params: LSCINT.P

Vial: 5  
Operator:  
Inst : Instrumen  
Multiplr: 1.00

Quant Method : C:\MSDCHEM\1...\HP022102.M (RTE Integrator)

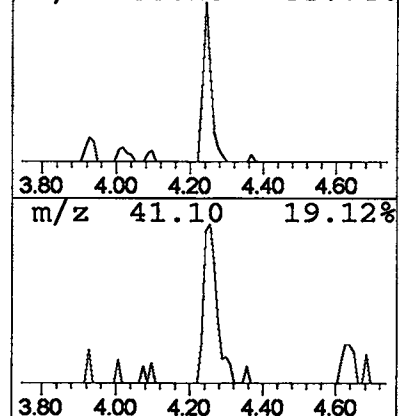
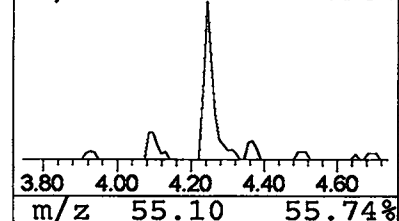
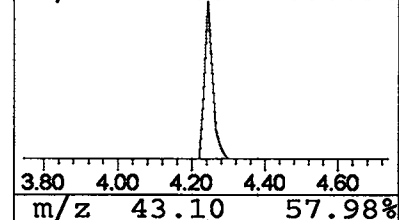
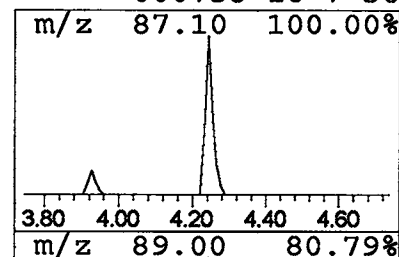
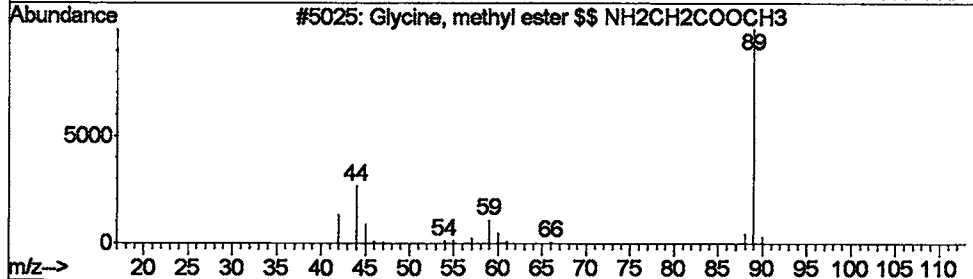
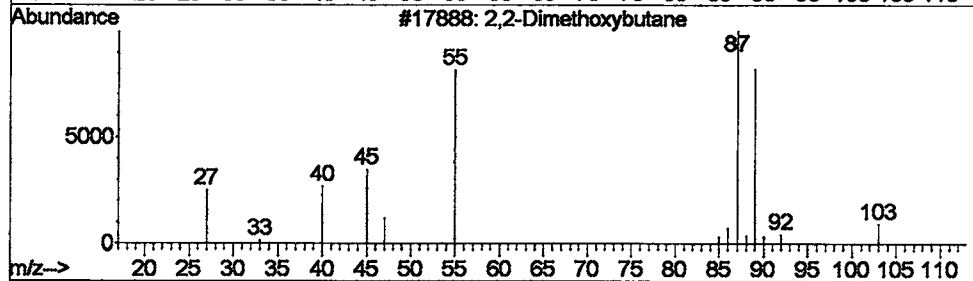
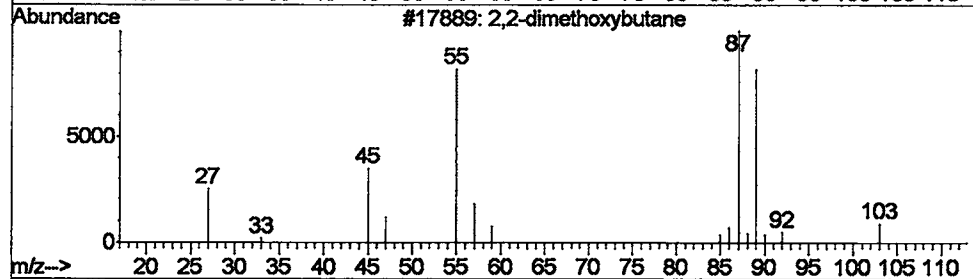
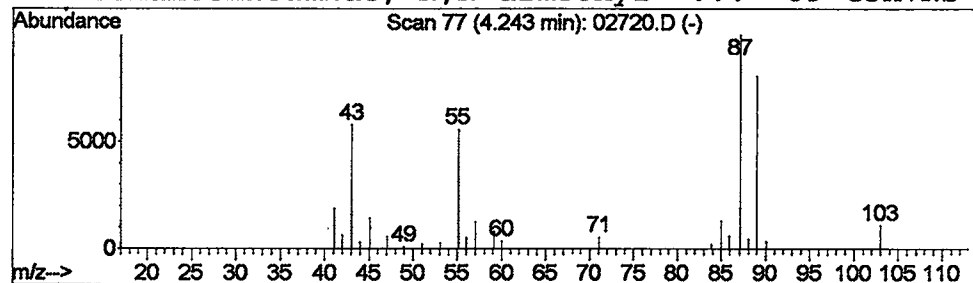
Title : 508 scan method

Library : C:\DATABASE\WILEY7N.L

\*\*\*\*\*  
Peak Number 3 2,2-dimethoxybutane Concentration Rank 3

| R.T. | EstConc   | Area   | Relative to ISTD       | R.T. |
|------|-----------|--------|------------------------|------|
| 4.24 | 0.44 ug/L | 188579 | 1,4-Dichlorobenzene-d4 | 9.71 |

| Hit# | of | Tentative ID                          | MW  | MolForm | CAS#        | Qual |
|------|----|---------------------------------------|-----|---------|-------------|------|
| 1    | 5  | 2,2-dimethoxybutane                   | 118 | C6H14O2 | 003453-99-4 | 83   |
| 2    |    | 2,2-Dimethoxybutane                   | 118 | C6H14O2 | 003453-99-4 | 74   |
| 3    |    | Glycine, methyl ester \$\$ NH2CH2C... | 89  | C3H7NO2 | 000616-34-2 | 45   |
| 4    |    | Methanethioamide, N,N-dimethyl- ...   | 89  | C3H7NS  | 000758-16-7 | 36   |



# Library Search Compound Report

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 Acq On : 28 Feb 2002 6:00 pm  
 Sample : 508 scan  
 Misc : February 28, 2002  
 MS Integration Params: LSCINT.P

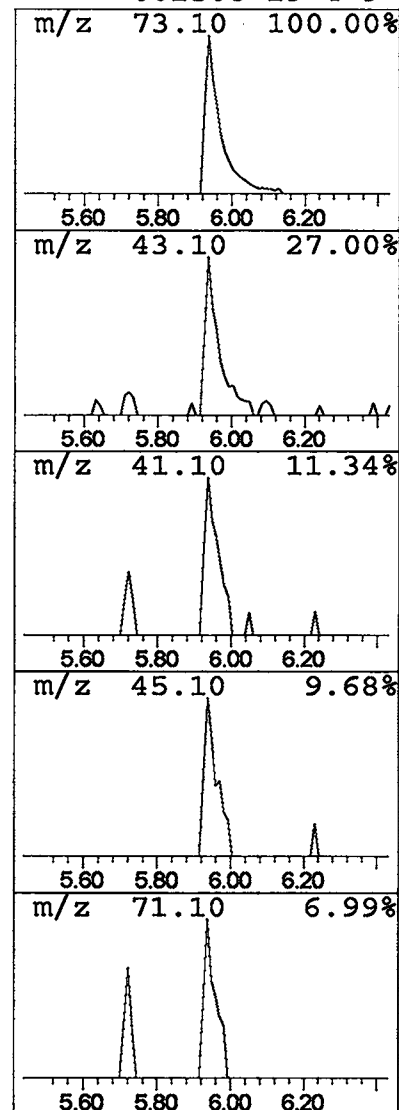
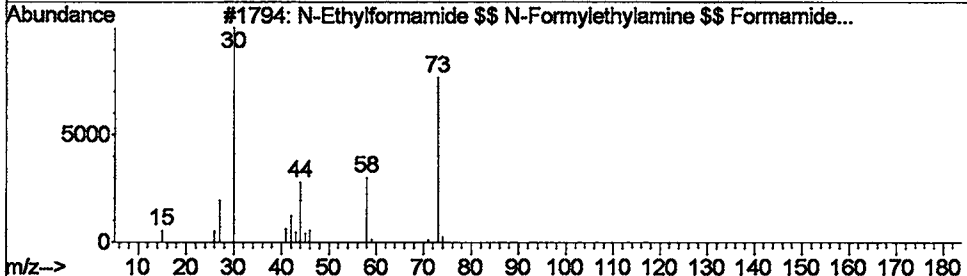
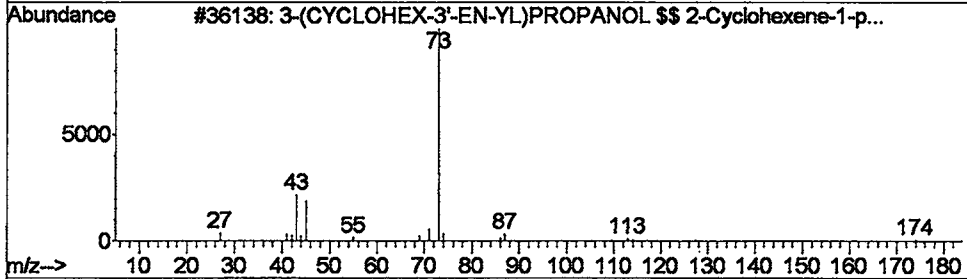
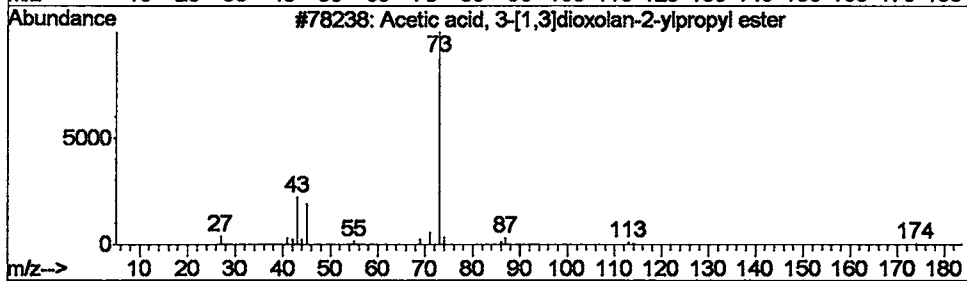
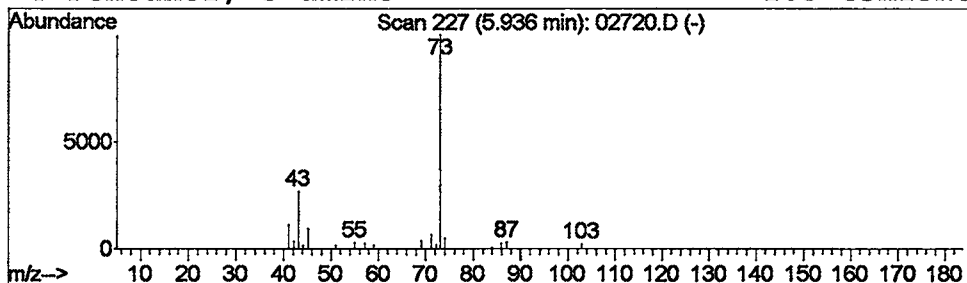
Vial: 5  
 Operator:  
 Inst : Instrumen  
 Multiplr: 1.00

Quant Method : C:\MSDCHEM\1...\HP022102.M (RTE Integrator)  
 Title : 508 scan method  
 Library : C:\DATABASE\WILEY7N.L

\*\*\*\*\*  
 Peak Number 4 Acetic acid, 3-[1,3]dioxola... Concentration Rank 4

| R.T. | EstConc   | Area   | Relative to ISTD       | R.T. |
|------|-----------|--------|------------------------|------|
| 5.94 | 0.38 ug/L | 160391 | 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       |   | N-Ethylformamide \$\$ N-Formylethy... | 73  | C3H7NO  | 000627-45-2 | 9    |
| 4       |   | Pentanol, 5-amino-                    | 103 | C5H13NO | 002508-29-4 | 9    |



# Library Search Compound Report

Data File : C:\MSDCHEM\1\5973N\02FEB28\02720.D  
 Acq On : 28 Feb 2002 6:00 pm  
 Sample : 508 scan  
 Misc : February 28, 2002  
 MS Integration Params: LSCINT.P

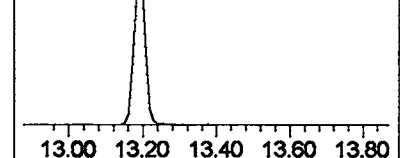
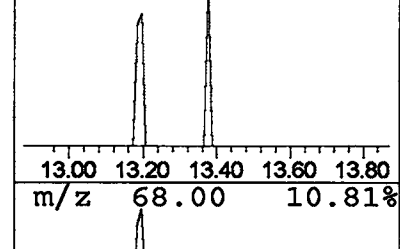
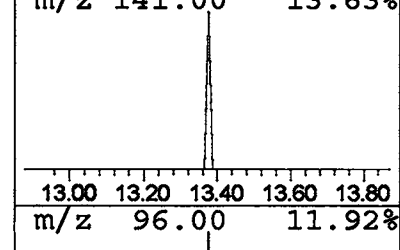
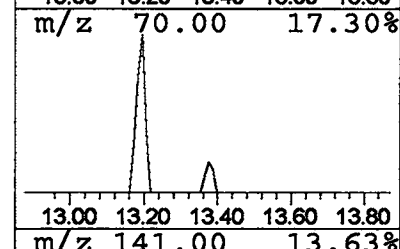
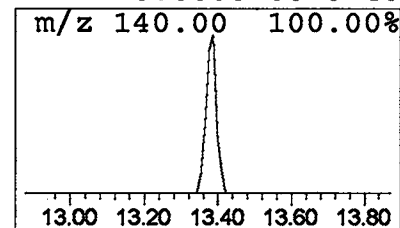
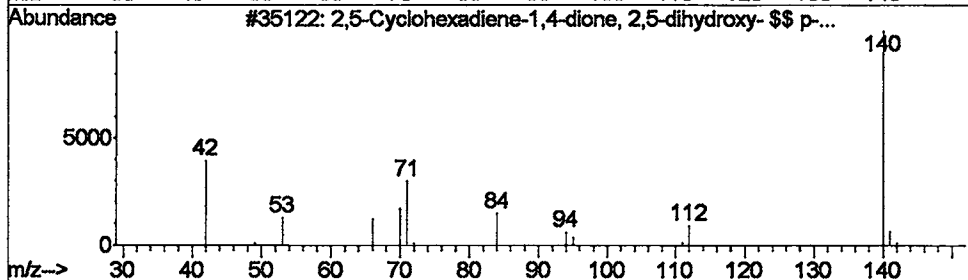
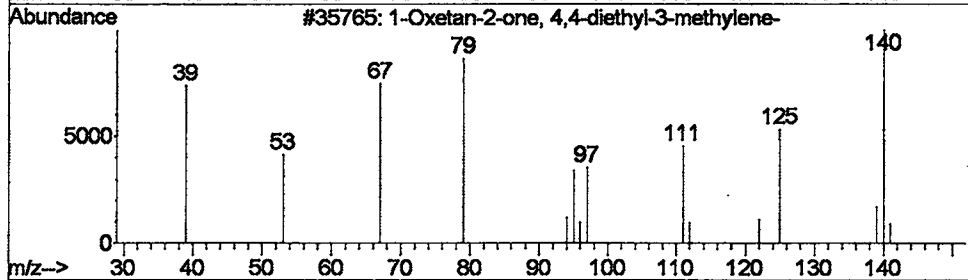
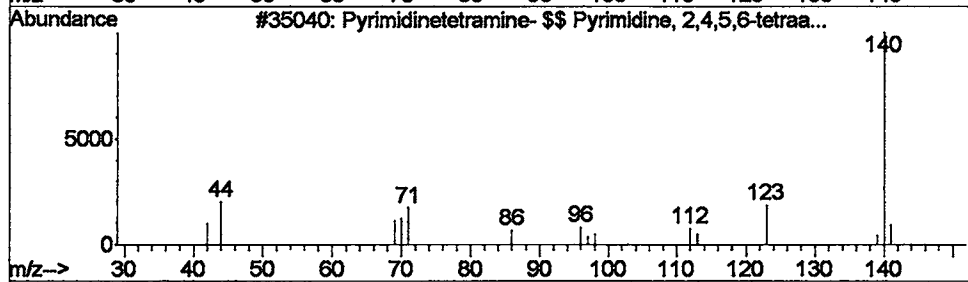
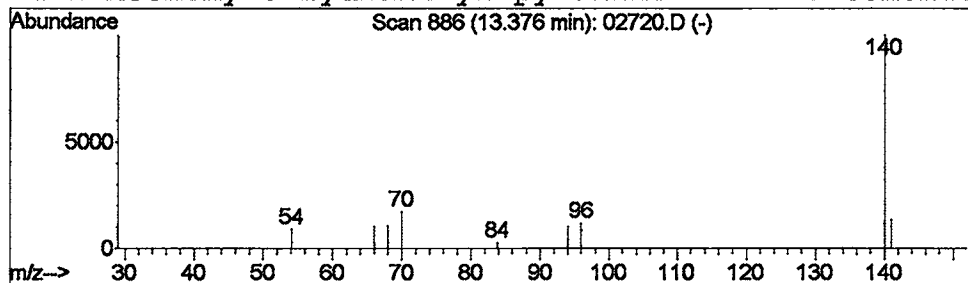
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 Operator:  
 Inst : Instrumen  
 Multiplr: 1.00

Quant Method : C:\MSDCHEM\1...\HP022102.M (RTE Integrator)  
 Title : 508 scan method  
 Library : C:\DATABASE\WILEY7N.L

\*\*\*\*\*  
 Peak Number 5 Pyrimidinetetramine- \$\$ Pyr... Concentration Rank 12

| R.T.  | EstConc   | Area  | Relative to ISTD | R.T.  |
|-------|-----------|-------|------------------|-------|
| 13.38 | 0.06 ug/L | 37307 | Naphthalene-d8   | 13.20 |

| Hit# | of | Tentative ID                          | MW  | MolForm | CAS#        | Qual |
|------|----|---------------------------------------|-----|---------|-------------|------|
| 1    | 5  | Pyrimidinetetramine- \$\$ Pyrimidi... | 140 | C4H8N6  | 001004-74-6 | 56   |
| 2    |    | 1-Oxetan-2-one, 4,4-diethyl-3-me...   | 140 | C8H12O2 | 000000-00-0 | 45   |
| 3    |    | 2,5-Cyclohexadiene-1,4-dione, 2,...   | 140 | C6H4O4  | 000615-94-1 | 40   |
| 4    |    | 2-Methoxy-3-hydrazinyl-pyrazine       | 140 | C5H8N4O | 000000-00-0 | 40   |



## Library Search Compound Report

Data File : C:\MSDCHEM\1\5973N\02FEB28\02720.D

Acq On : 28 Feb 2002 6:00 pm

Sample : 508 scan

Misc : February 28, 2002

MS Integration Params: LSCINT.P

Vial: 5

Operator:

Inst : Instrumen

Multiplr: 1.00

Quant Method : C:\MSDCHEM\1...\HP022102.M (RTE Integrator)

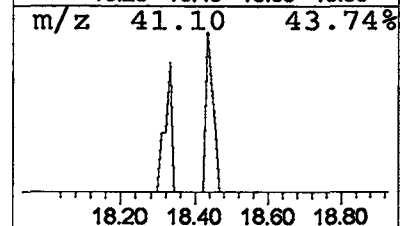
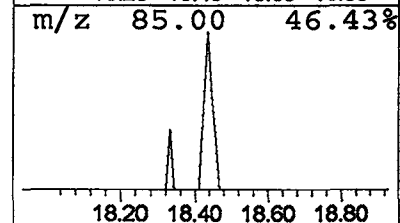
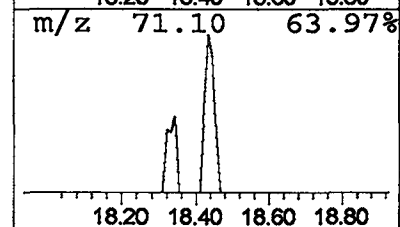
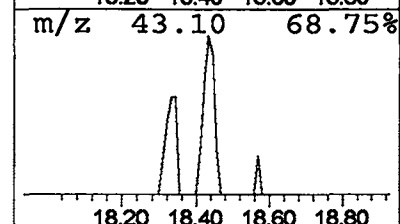
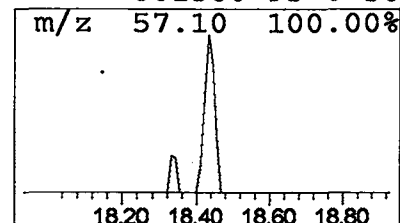
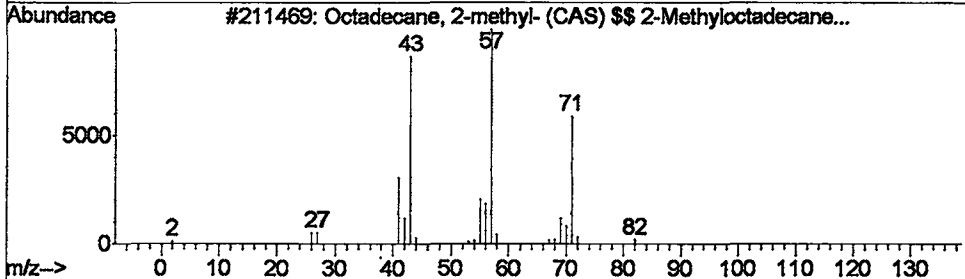
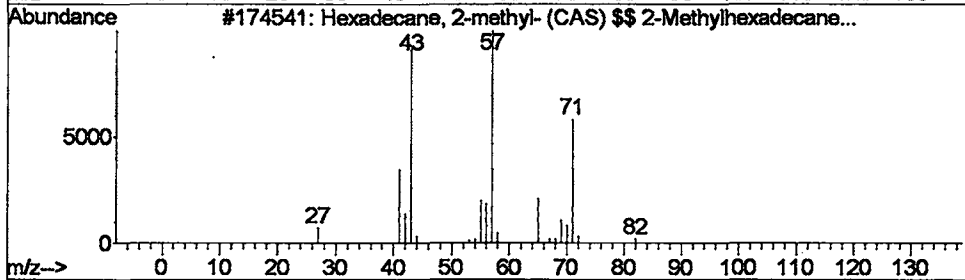
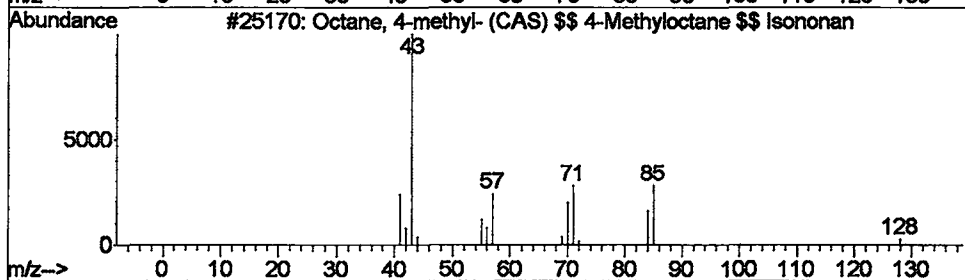
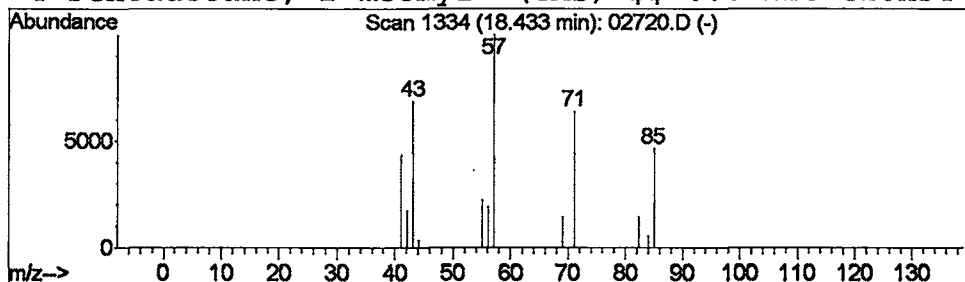
Title : 508 scan method

Library : C:\DATABASE\WILEY7N.L

\*\*\*\*\*  
Peak Number 6 Octane, 4-methyl- (CAS) \$\$ ... Concentration Rank 9

| R.T.  | EstConc   | Area  | Relative to ISTD | R.T.  |
|-------|-----------|-------|------------------|-------|
| 18.43 | 0.10 ug/L | 66680 | Acenaphthene-d10 | 18.34 |

| Hit# of 5 | Tentative ID                          | MW  | MolForm | CAS#        | Qual |
|-----------|---------------------------------------|-----|---------|-------------|------|
| 1         | Octane, 4-methyl- (CAS) \$\$ 4-Met... | 128 | C9H20   | 002216-34-4 | 53   |
| 2         | Hexadecane, 2-methyl- (CAS) \$\$ 2... | 240 | C17H36  | 001560-92-5 | 50   |
| 3         | Octadecane, 2-methyl- (CAS) \$\$ 2... | 268 | C19H40  | 001560-88-9 | 50   |
| 4         | Pentadecane, 2-methyl- (CAS) \$\$ ... | 226 | C16H34  | 001560-93-6 | 50   |



# Library Search Compound Report

Data File : C:\MSDCHEM\1\5973N\02FEB28\02720.D  
 Acq On : 28 Feb 2002 6:00 pm  
 Sample : 508 scan  
 Misc : February 28, 2002  
 MS Integration Params: LSCINT.P

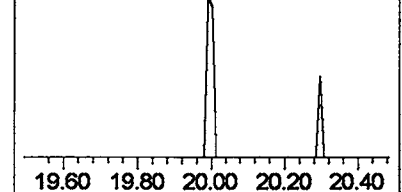
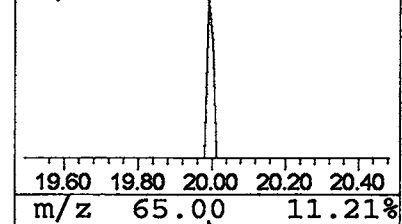
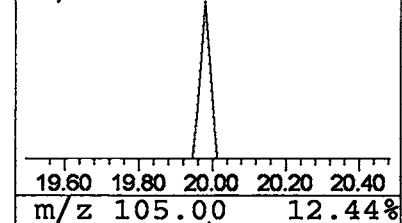
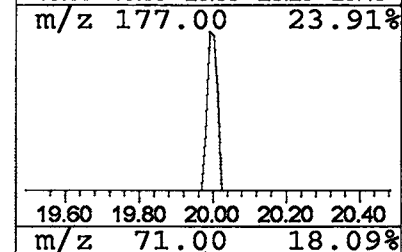
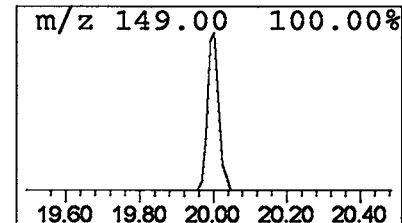
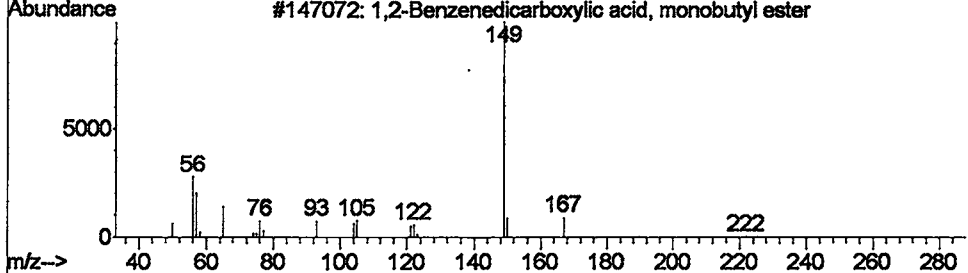
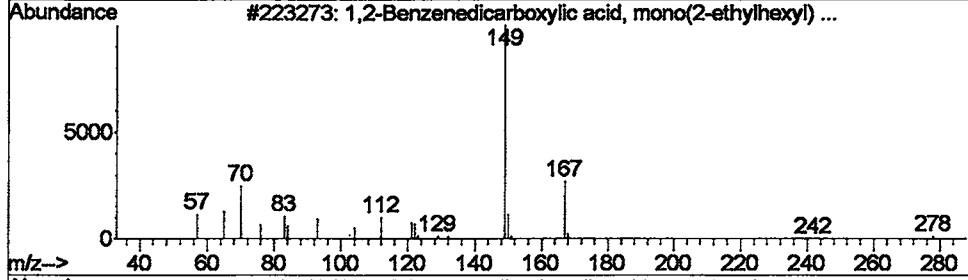
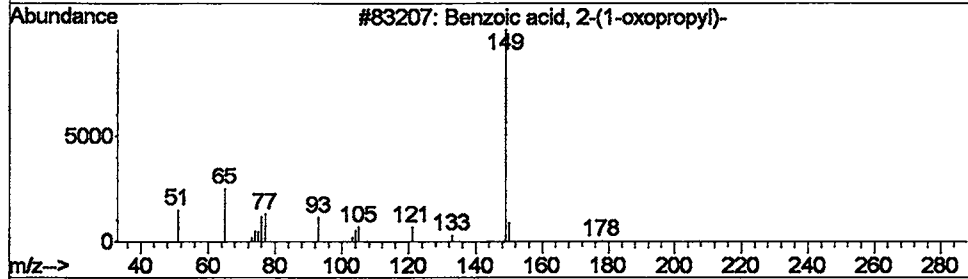
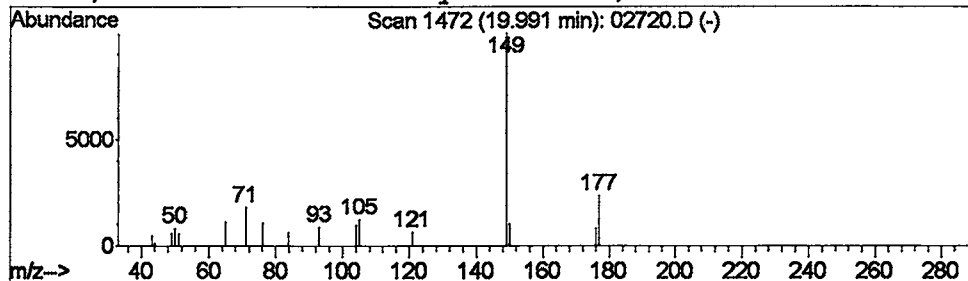
Vial: 5  
 Operator:  
 Inst : Instrumen  
 Multiplr: 1.00

Quant Method : C:\MSDCHEM\1...\HP022102.M (RTE Integrator)  
 Title : 508 scan method  
 Library : C:\DATABASE\WILEY7N.L

\*\*\*\*\*  
 Peak Number 7 Benzoic acid, 2-(1-oxopropyl)- Concentration Rank 10

| R.T.  | EstConc   | Area  | Relative to ISTD | R.T.  |
|-------|-----------|-------|------------------|-------|
| 19.99 | 0.08 ug/L | 52230 | Acenaphthene-d10 | 18.34 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm  | CAS#        | Qual |
|------|----|---|-------------------------------------|-----|----------|-------------|------|
| 1    |    |   | Benzoic acid, 2-(1-oxopropyl)-      | 178 | C10H10O3 | 002360-45-4 | 64   |
| 2    |    |   | 1,2-Benzenedicarboxylic acid, mo... | 278 | C16H22O4 | 004376-20-9 | 59   |
| 3    |    |   | 1,2-Benzenedicarboxylic acid, mo... | 222 | C12H14O4 | 000131-70-4 | 59   |
| 4    |    |   | 1,2-Benzenedicarboxylic acid, mo... | 222 | C12H14O4 | 000131-70-4 | 59   |



# Library Search Compound Report

Data File : C:\MSDCHEM\1\5973N\02FEB28\02720.D  
 Acq On : 28 Feb 2002 6:00 pm  
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 Misc : February 28, 2002  
 MS Integration Params: LSCINT.P

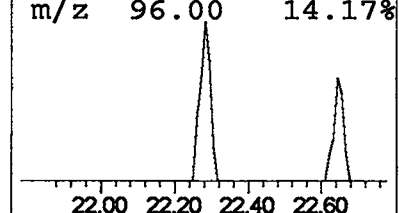
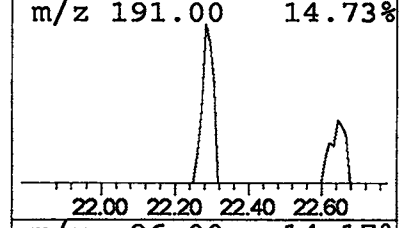
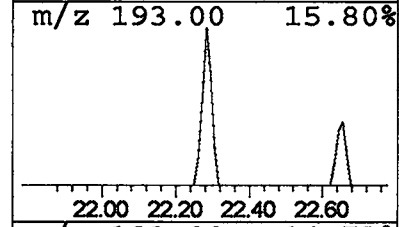
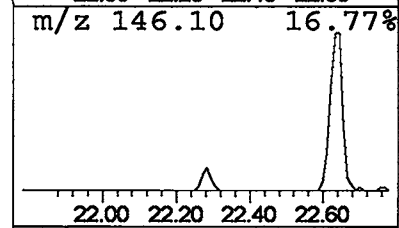
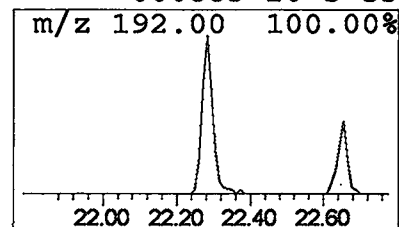
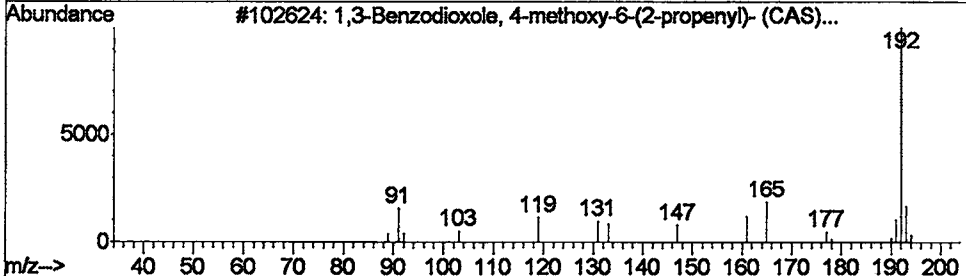
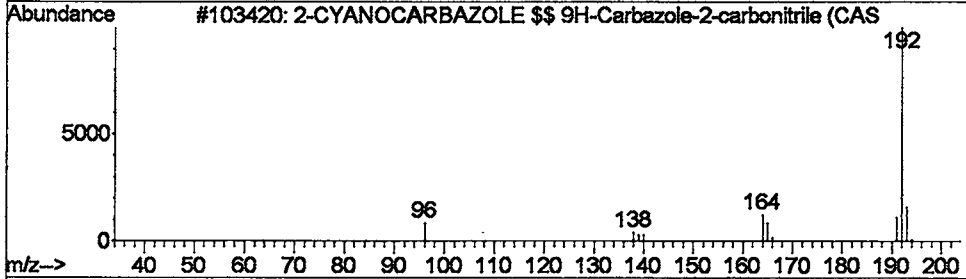
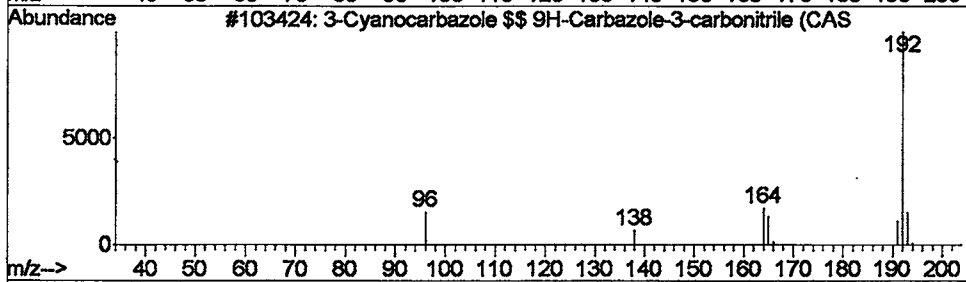
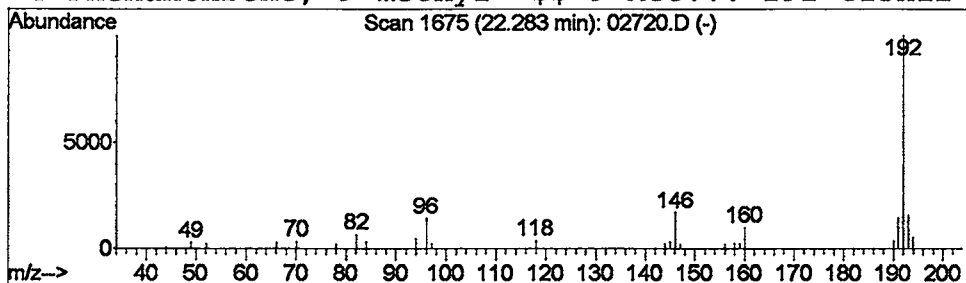
Vial: 5  
 Operator:  
 Inst : Instrumen  
 Multiplr: 1.00

Quant Method : C:\MSDCHEM\1...\HP022102.M (RTE Integrator)  
 Title : 508 scan method  
 Library : C:\DATABASE\WILEY7N.L

\*\*\*\*\*  
 Peak Number 8 3-Cyanocarbazole \$\$ 9H-Carb... Concentration Rank 7

| R.T.  | EstConc   | Area   | Relative to ISTD | R.T.  |
|-------|-----------|--------|------------------|-------|
| 22.28 | 0.14 ug/L | 101679 | Phenanthrene-d10 | 22.63 |

| Hit# of | 5                                     | Tentative ID | MW       | MolForm     | CAS# | Qual |
|---------|---------------------------------------|--------------|----------|-------------|------|------|
| 1       | 3-Cyanocarbazole \$\$ 9H-Carbazole... | 192          | C13H8N2  | 057102-93-9 | 64   |      |
| 2       | 2-CYANOCARBAZOLE \$\$ 9H-Carbazole... | 192          | C13H8N2  | 057955-18-7 | 64   |      |
| 3       | 1,3-Benzodioxole, 4-methoxy-6-(2...   | 192          | C11H12O3 | 000607-91-0 | 59   |      |
| 4       | Phenanthrene, 9-methyl- \$\$ 9-Met... | 192          | C15H12   | 000883-20-5 | 53   |      |



# Library Search Compound Report

Data File : C:\MSDCHEM\1\5973N\02FEB28\02720.D  
 Acq On : 28 Feb 2002 6:00 pm  
 Sample : 508 scan  
 Misc : February 28, 2002  
 MS Integration Params: LSCINT.P

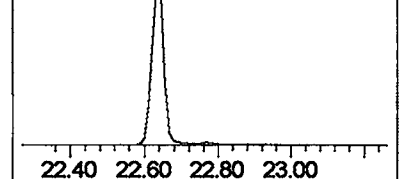
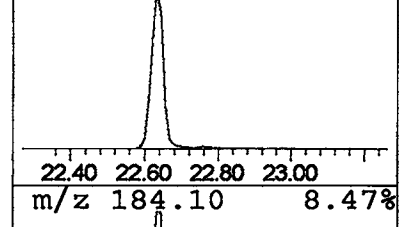
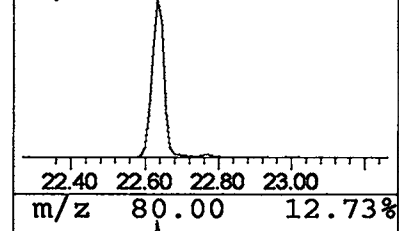
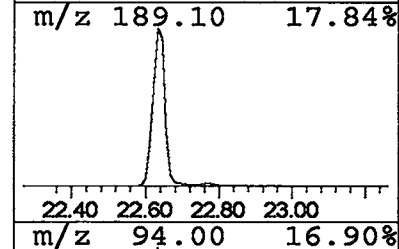
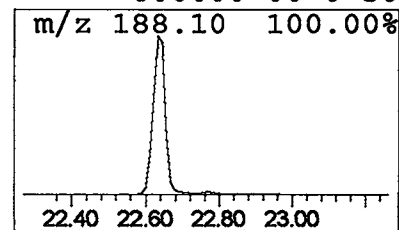
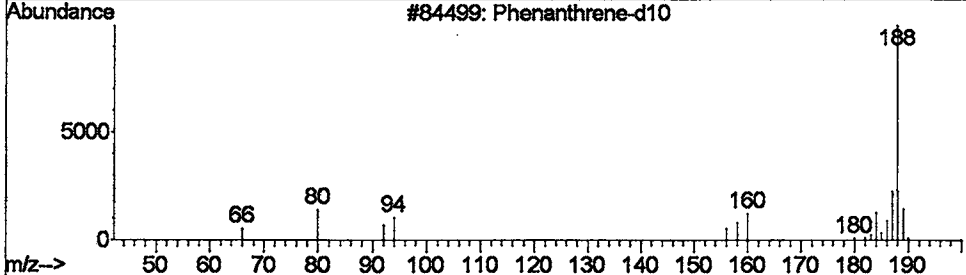
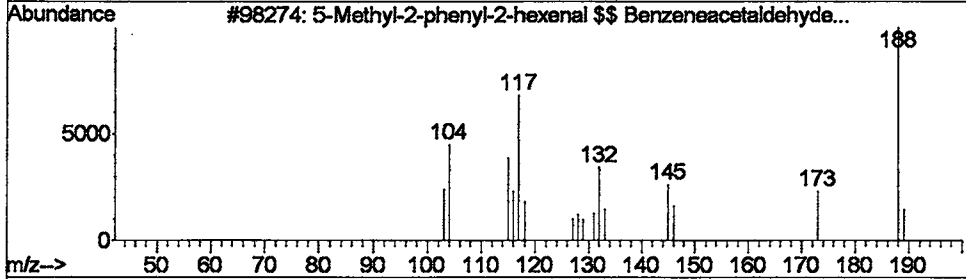
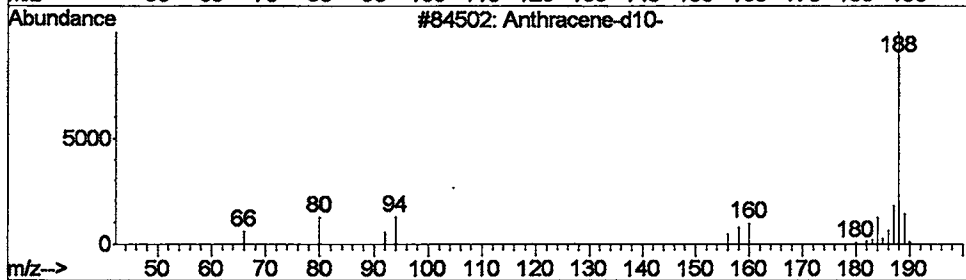
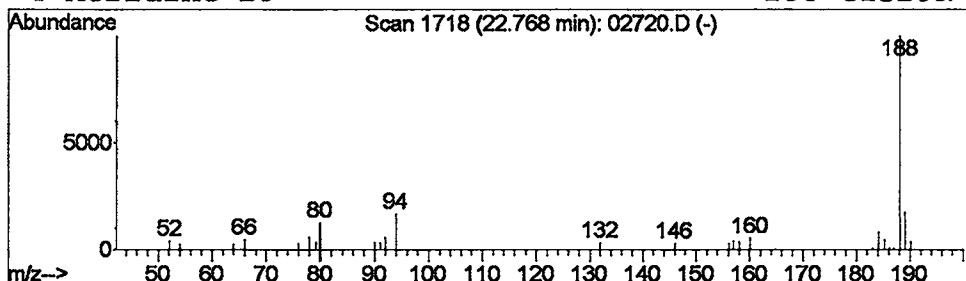
Vial: 5  
 Operator:  
 Inst : Instrumen  
 Multiplr: 1.00

Quant Method : C:\MSDCHEM\1...\HP022102.M (RTE Integrator)  
 Title : 508 scan method  
 Library : C:\DATABASE\WILEY7N.L

\*\*\*\*\*  
 Peak Number 9 Anthracene-d10- Concentration Rank 2

| R.T.  | EstConc   | Area   | Relative to ISTD | R.T.  |
|-------|-----------|--------|------------------|-------|
| 22.77 | 0.53 ug/L | 387292 | Phenanthrene-d10 | 22.63 |

| Hit# | of | Tentative ID                          | MW  | MolForm | CAS#        | Qual |
|------|----|---------------------------------------|-----|---------|-------------|------|
| 1    | 5  | Anthracene-d10-                       | 188 | C14D10  | 001719-06-8 | 64   |
| 2    |    | 5-Methyl-2-phenyl-2-hexenal \$\$ B... | 188 | C13H16O | 021834-92-4 | 52   |
| 3    |    | Phenanthrene-d10                      | 188 | C14D10  | 001517-22-2 | 50   |
| 4    |    | Acridine-D9                           | 188 | C13D9N  | 000000-00-0 | 50   |



# Library Search Compound Report

Data File : C:\MSDCHEM\1\5973N\02FEB28\02720.D  
 Acq On : 28 Feb 2002 6:00 pm  
 Sample : 508 scan  
 Misc : February 28, 2002  
 MS Integration Params: LSCINT.P

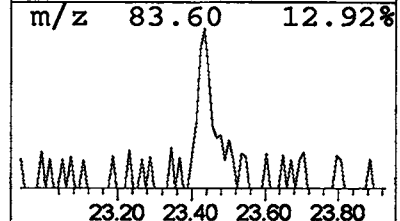
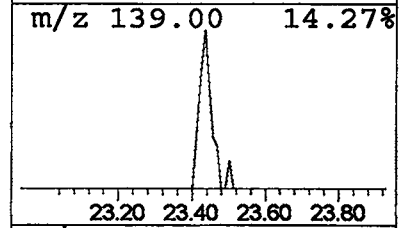
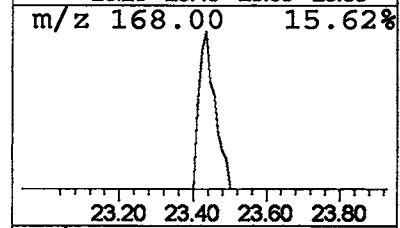
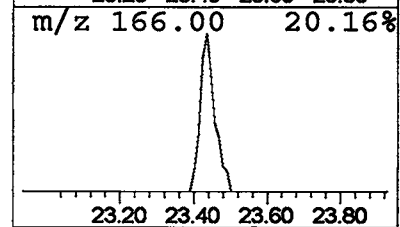
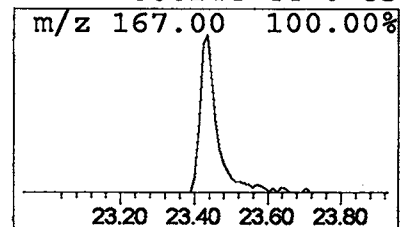
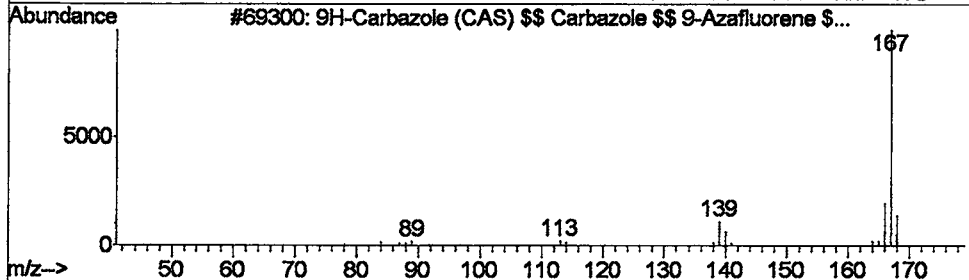
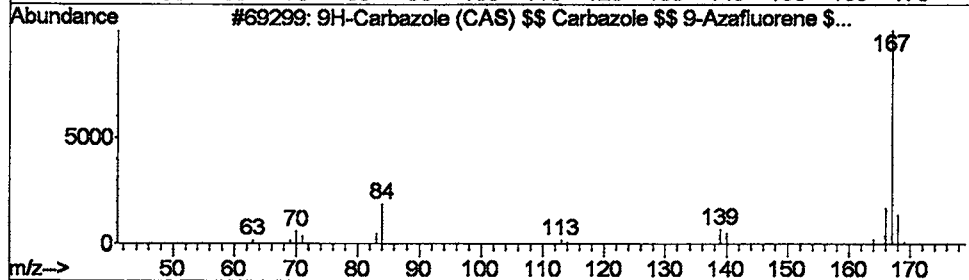
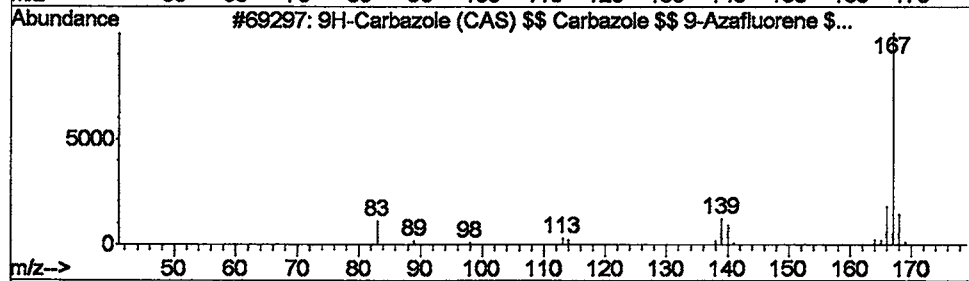
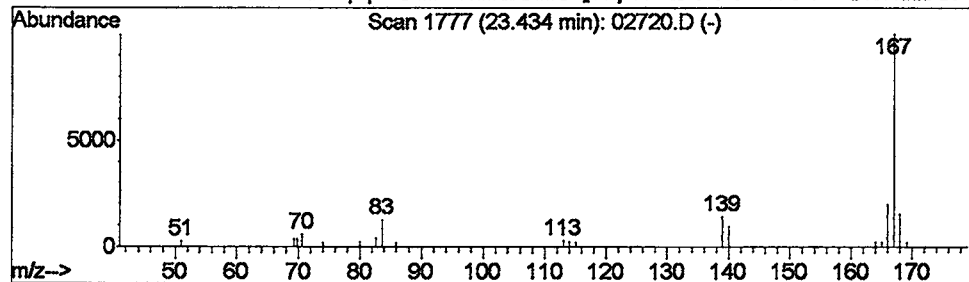
Vial: 5  
 Operator:  
 Inst : Instrumen  
 Multiplr: 1.00

Quant Method : C:\MSDCHEM\1...\HP022102.M (RTE Integrator)  
 Title : 508 scan method  
 Library : C:\DATABASE\WILEY7N.L

\*\*\*\*\*  
 Peak Number 10 9H-Carbazole (CAS) \$\$ Carba... Concentration Rank 5

| R.T.  | EstConc   | Area   | Relative to ISTD | R.T.  |
|-------|-----------|--------|------------------|-------|
| 23.43 | 0.18 ug/L | 129397 | Phenanthrene-d10 | 22.63 |

| Hit# | of | Tentative ID                          | MW  | MolForm | CAS#        | Qual |
|------|----|---------------------------------------|-----|---------|-------------|------|
| 1    | 5  | 9H-Carbazole (CAS) \$\$ Carbazole ... | 167 | C12H9N  | 000086-74-8 | 90   |
| 2    |    | 9H-Carbazole (CAS) \$\$ Carbazole ... | 167 | C12H9N  | 000086-74-8 | 86   |
| 3    |    | 9H-Carbazole (CAS) \$\$ Carbazole ... | 167 | C12H9N  | 000086-74-8 | 83   |
| 4    |    | 1-Azafluorene \$\$ 9H-Indeno[2,1-b... | 167 | C12H9N  | 000244-44-0 | 83   |



## Library Search Compound Report

Data File : C:\MSDCHEM\1\5973N\02FEB28\02720.D  
Acq On : 28 Feb 2002 6:00 pm  
Sample : 508 scan  
Misc : February 28, 2002  
MS Integration Params: LSCINT.P

Vial: 5  
Operator:  
Inst : Instrumen  
Multiplr: 1.00

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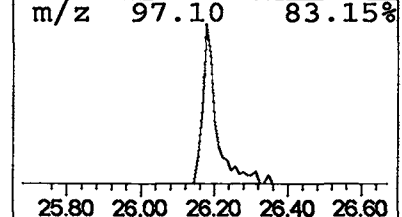
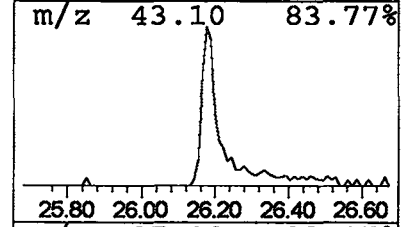
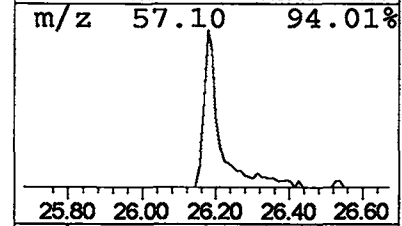
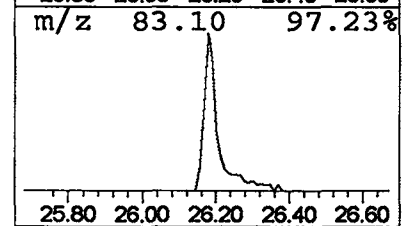
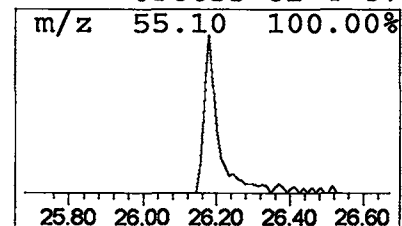
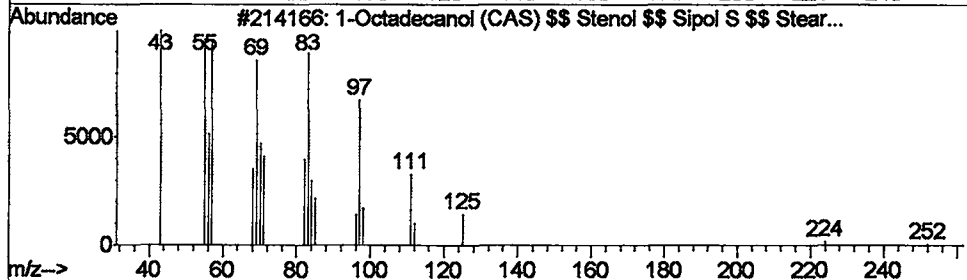
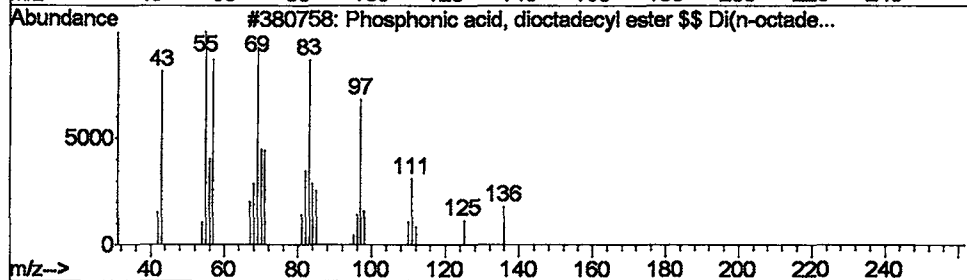
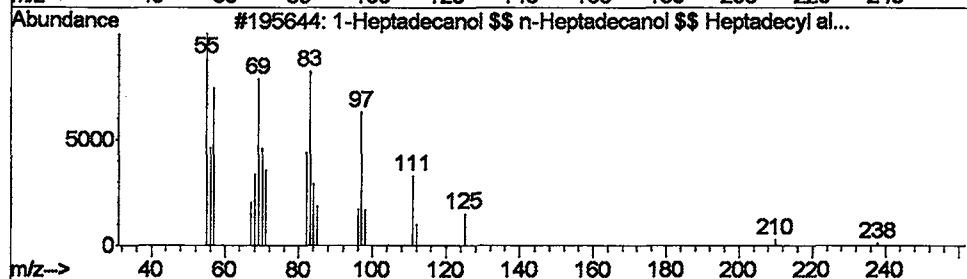
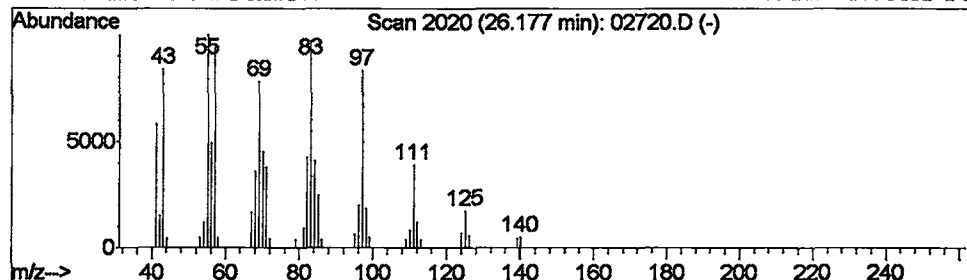
Title : 508 scan method

Library : C:\DATABASE\WILEY7N.L

\*\*\*\*\*  
Peak Number 11 1-Heptadecanol \$\$ n-Heptade... Concentration Rank 1

| R.T.  | EstConc   | Area   | Relative to ISTD | R.T.  |
|-------|-----------|--------|------------------|-------|
| 26.18 | 0.75 ug/L | 541745 | Phenanthrene-d10 | 22.63 |

| Hit# | of | 5 | Tentative ID                            | MW  | MolForm   | CAS#        | Qual |
|------|----|---|-----------------------------------------|-----|-----------|-------------|------|
| 1    |    |   | 1-Heptadecanol \$\$ n-Heptadecanol...   | 256 | C17H36O   | 001454-85-9 | 94   |
| 2    |    |   | Phosphonic acid, dioctadecyl est...     | 587 | C36H75O3P | 019047-85-9 | 93   |
| 3    |    |   | 1-Octadecanol (CAS) \$\$ Stenol \$\$... | 270 | C18H38O   | 000112-92-5 | 90   |
| 4    |    |   | 1-Hexadecanol                           | 242 | C16H34O   | 036653-82-4 | 87   |



## Tentatively Identified Compound (LSC) summary

Operator ID:            Date Acquired: 28 Feb 2002    6:00 pm  
 Data File: C:\MSDCHEM\1\5973N\02FEB28\02720.D  
 Name: 508 scan  
 Misc: February 28, 2002  
 Method: C:\MSDCHEM\1\METHODS\METHODS\HP022102.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...              | 3.62  | 0.2     | ug/L  | 65579    | 1                     | 9.71  | 8513120  | 20.0 |
| Propane, 1,1-dime...              | 3.93  | 0.1     | ug/L  | 29091    | 1                     | 9.71  | 8513120  | 20.0 |
| 2,2-dimethoxybutane               | 4.24  | 0.4     | ug/L  | 188579   | 1                     | 9.71  | 8513120  | 20.0 |
| Acetic acid, 3-[1...              | 5.94  | 0.4     | ug/L  | 160391   | 1                     | 9.71  | 8513120  | 20.0 |
| Pyrimidinetetrami...              | 13.38 | 0.1     | ug/L  | 37307    | 2                     | 13.20 | 12479800 | 20.0 |
| Octane, 4-methyl-...              | 18.43 | 0.1     | ug/L  | 66680    | 3                     | 18.34 | 13588300 | 20.0 |
| Benzoic acid, 2-(...              | 19.99 | 0.1     | ug/L  | 52230    | 3                     | 18.34 | 13588300 | 20.0 |
| 3-Cyanocarbazole ...              | 22.28 | 0.1     | ug/L  | 101679   | 4                     | 22.63 | 14490100 | 20.0 |
| <del>Anthracene-d10-</del>        | 22.77 | 0.5     | ug/L  | 387292   | 4                     | 22.63 | 14490100 | 20.0 |
| <del>9H-Carbazole (CAS...</del>   | 23.43 | 0.2     | ug/L  | 129397   | 4                     | 22.63 | 14490100 | 20.0 |
| <del>1-Heptadecanol \$\$...</del> | 26.18 | 0.7     | ug/L  | 541745   | 4                     | 22.63 | 14490100 | 20.0 |
| Pyrene (CAS) \$\$ ....            | 26.84 | 0.1     | ug/L  | 96136    | 5                     | 30.49 | 13760200 | 20.0 |

Westchester  
gov.comENVIRONMENTAL SERVICES SAMPLE SUBMISSION FORM  
WESTCHESTER COUNTY LABORATORIES AND RESEARCH2 Dana Road, Valhalla, New York 10595  
(914) 231-1620 Fax (914) 231-1772NYS-ELAP No. 10108  
Form A-70 (rev.10/01)

Time/Date Set:

## Sample Type (circle one):

1. Potable 3. Non Potable Treated  
2. Non Potable 4. Other \_\_\_\_\_

Time/Date Collected: 1230 2-11-02

Collected By: [Signature] Agency Code: \_\_\_\_\_

## Collected From:

Location's/Individual's Name: LYO - BCP

Street: 19 W. LAKE DR.

City/St/Zip: Valhalla ny 10595

Identification of Source: \_\_\_\_\_

Collection Point: CLO 6 H

UOA, 1487, 13, 14, 15 C-1480  
 22574, 2570, 2545, 2516, 2597, 2512, 2597  
 Bottle #s: P-1482, 2750, 2764

Chain of Custody? Yes No

## Comments:

PWS #: \_\_\_\_\_ Source ID: \_\_\_\_\_ Type Descriptor: \_\_\_\_\_

## Bill to:

Name/Company: \_\_\_\_\_

Street: \_\_\_\_\_

City/St/Zip: \_\_\_\_\_

Phone#: \_\_\_\_\_ Fax#: \_\_\_\_\_

CA,CH, BI? \_\_\_\_\_ Check#: \_\_\_\_\_ Amount\$ \_\_\_\_\_

## BACTERIOLOGICAL ANALYSIS

- Lab#: \_\_\_\_\_
- \_\_\_ Heterotrophic Plate Count (hemodialysis & all others)  
 \_\_\_ Coliform P/A (Public supplies)  
 \_\_\_ Coliform MPN (priv well, raw, stp)  
 \_\_\_ Fecal Coliform (all samples)  
 \_\_\_ Microscopic  
 \_\_\_ Macroscopic  
 \_\_\_ API  
 \_\_\_ Other: \_\_\_\_\_

Refrigerated? yes no

Chlorinated: yes no

Free Chlorine Res: \_\_\_\_\_ mg/L

Total Chlorine Res: \_\_\_\_\_ mg/L

pH: \_\_\_\_\_

## RADIOCHEMICAL ANALYSIS

- \_\_\_ Radon  
 \_\_\_ Other: \_\_\_\_\_

Lab#: \_\_\_\_\_

## ORGANIC POTABLE ANALYSIS

- \_\_\_ Trihalomethanes (524)  
☒ Volatile Halocarbons (524)  
☒ Volatile Aromatics (524)  
 \_\_\_ Total Trihalomethane Potential (510)  
 \_\_\_ Microextractables (504)  
 \_\_\_ Pesticides/PCB screen (508)  
 \_\_\_ PCB as Decachlorobiphenyl (508A)  
 \_\_\_ Herbicides (515)  
 \_\_\_ Pesticides (525)  
 \_\_\_ Carbamate Pesticides (531)  
 \_\_\_ Glyphosate (547)  
 \_\_\_ Diquat (549)  
 \_\_\_ Chlorination Disinfection Byproducts (551)  
 \_\_\_ Haloacetic Acids (552)  
 \_\_\_ Other Pesticides (SM18/6630)  
 \_\_\_ UCMR List 1  
 \_\_\_ MtBE Only (524)

## ORGANIC NONPOTABLE/SOLID ANALYSIS

- \_\_\_ Purgeable Halocarbons (624/8260)  
 \_\_\_ Purgeable Aromatics (624/8260)  
 \_\_\_ Acrolein & Acrylonitrile (8260)  
 \_\_\_ Organochlorine Pesticides (608/8081)  
 \_\_\_ PCB's (Arochors) (608/8081)  
 \_\_\_ Herbicides (SM18/6630,8151)  
 \_\_\_ Other Pesticides (SM18/6630)  
 \_\_\_ Petroleum Hydrocarbon Scan (310-13)  
 \_\_\_ Glycols in Water  
 \_\_\_ Acid Extractables (625/8270)  
 \_\_\_ Base Neutral Extractables (625/8270)  
☒ GC/MS Unknown Scan  
 \_\_\_ Phthalates (606MS)  
 \_\_\_ Other: \_\_\_\_\_

## INORGANIC ANALYSIS REQUESTED - Lab#: 2720

- \_\_\_ Color  
 \_\_\_ Turbidity  
 \_\_\_ pH  
 \_\_\_ Conductivity  
 \_\_\_ Corrosivity (temp: \_\_\_\_\_ °C)  
 \_\_\_ BOD (5 day)  
 \_\_\_ Carbonaceous BOD (5 day)  
 \_\_\_ Soluble BOD (5 day)  
 \_\_\_ Chemical Oxygen Demand  
 \_\_\_ Dissolved Oxygen  
 \_\_\_ Total Organic Carbon  
 \_\_\_ Total Organic Halides  
 \_\_\_ Solids, Percent (%)  
 \_\_\_ Solids, Settleable  
 \_\_\_ Solids, Suspended  
 \_\_\_ Solids, Total  
 \_\_\_ Solids, Total Dissolved  
 \_\_\_ Solids, Total Volatile  
☒ Other: 7A

- \_\_\_ Acidity  
 \_\_\_ Alkalinity  
 \_\_\_ Bromide  
 \_\_\_ Bromate/Chlorate/Chlorite  
 \_\_\_ Chloride  
 \_\_\_ Cyanide  
 \_\_\_ Fluoride  
 \_\_\_ Hardness, Total as CaCO<sub>3</sub>  
 \_\_\_ Nitrogen, Ammonia as N  
 \_\_\_ Nitrogen, Nitrate as N  
 \_\_\_ Nitrogen, Nitrite as N  
 \_\_\_ Nitrogen, Total Kjeldahl as N  
 \_\_\_ Oil and Grease  
 \_\_\_ Phenol  
 \_\_\_ Phosphorus Total as P  
 \_\_\_ Phosphorus, Ortho as P  
 \_\_\_ Sulfates  
 \_\_\_ Surfactants  
 \_\_\_ UV254

- ☒ Aluminum  
☒ Antimony  
☒ Arsenic  
☒ Barium  
☒ Beryllium  
☒ Boron  
☒ Cadmium  
 \_\_\_ Calcium, as CaCO<sub>3</sub>  
 \_\_\_ Calcium, Ca<sup>+2</sup>  
☒ Chromium, Total  
☒ Chromium, Hexavalent  
☒ Cobalt  
☒ Copper  
☒ Iron  
☒ Lead  
 \_\_\_ Magnesium  
☒ Manganese  
☒ Mercury  
☒ Molybdenum  
☒ Nickel

- \_\_\_ Potassium  
☒ Selenium  
☒ Silver  
 \_\_\_ Sodium  
☒ Thallium  
 \_\_\_ Vanadium  
☒ Zinc

## CLP

- \_\_\_ Metals/CN  
 \_\_\_ Volatiles  
 \_\_\_ Semi-volatiles  
 \_\_\_ Pesticides/PCB's

## TCLP

- \_\_\_ TCLP-Metals  
 \_\_\_ TCLP-Pesticides  
 \_\_\_ TCLP-Herbicides  
 \_\_\_ TCLP-Volatiles  
 \_\_\_ TCLP-BNA's