

User: Vinci, David L  
Westchester Dept. of Labs & Research  
Date: 05/20/2002 Time: 13:59:46

Sample: AE10704

Analysis: \$GCMS GCMS Scan For Unknowns

Units: ug

Started: 5/20/02 13:59 Ended: 5/20/02 13:59 Analyst: DLV

Page: 1

|   | Component Name          | Viol | Result | Qualifier | MDL  |
|---|-------------------------|------|--------|-----------|------|
| 1 | Other unknown compounds |      | .      |           |      |
| 2 | Surr_2,4-DDD            |      | 75     |           | 10.0 |

**Comments for Sample AE10704 - Analysis \$GCMS**

Westchester Dept. of Labs & Research

User: Vinci, David L

Date: 05-20-2002 Time: 14:00:25

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

Data File : C:\MSDCHEM\1\DATA\020513\10704.D

Vial: 15

Acq On : 14 May 2002 2:23 am

Operator: Nefliu

Sample : sample

Inst : Instrumen

Misc :

Multiplr: 1.00

MS Integration Params: rteint.e

Quant Time: May 15 08:59:28 2002

Quant Results File: SCAN020507.RE

Quant Method : C:\MSDCHEM\1\METHODS\SCAN020507.M (RTE Integrator)

Title : Method 508 GC/MS Scan

Last Update : Thu May 09 09:21:01 2002

Response via : Initial Calibration

DataAcq Meth : 508SCAN1

| Internal Standards        | R.T.  | QIon | Response | Conc  | Units | Dev(Min) |
|---------------------------|-------|------|----------|-------|-------|----------|
| 1) 1,4-Dichlorobenzene-d4 | 11.38 | 152  | 554984   | 40.00 | ug/l  | -0.01    |
| 10) Naphthalene-d8        | 16.54 | 136  | 2091779  | 40.00 | ug/l  | -0.03    |
| 17) Acenaphthene-d10      | 24.47 | 164  | 1123816  | 40.00 | ug/l  | -0.01    |
| 30) Phenanthrene-d10      | 31.72 | 188  | 1905232  | 40.00 | ug/l  | -0.02    |
| 38) Chrysene-d12          | 39.94 | 240  | 1399208  | 40.00 | ug/l  | -0.05    |
| 44) Perylene-d12          | 43.16 | 264  | 864902   | 40.00 | ug/l  | -0.02    |

System Monitoring Compounds

|               |        |     |          |      |        |       |
|---------------|--------|-----|----------|------|--------|-------|
| 27) TCMX      | 27.54  | 207 | 67204    | 7.49 | ug/l   | -0.04 |
| Spiked Amount | 10.000 |     | Recovery | =    | 74.90% |       |

Target Compounds

|                                |       |     |       |      |      |              |
|--------------------------------|-------|-----|-------|------|------|--------------|
| 43) Bis(2-ethylhexyl)phthalate | 40.52 | 149 | 15063 | 0.50 | ug/l | Qvalue<br>94 |
|--------------------------------|-------|-----|-------|------|------|--------------|

-----  
 (#) = qualifier out of range (m) = manual integration (+) = signals summed  
 10704.D SCAN020507.M Wed May 15 09:01:03 2002 Page 1

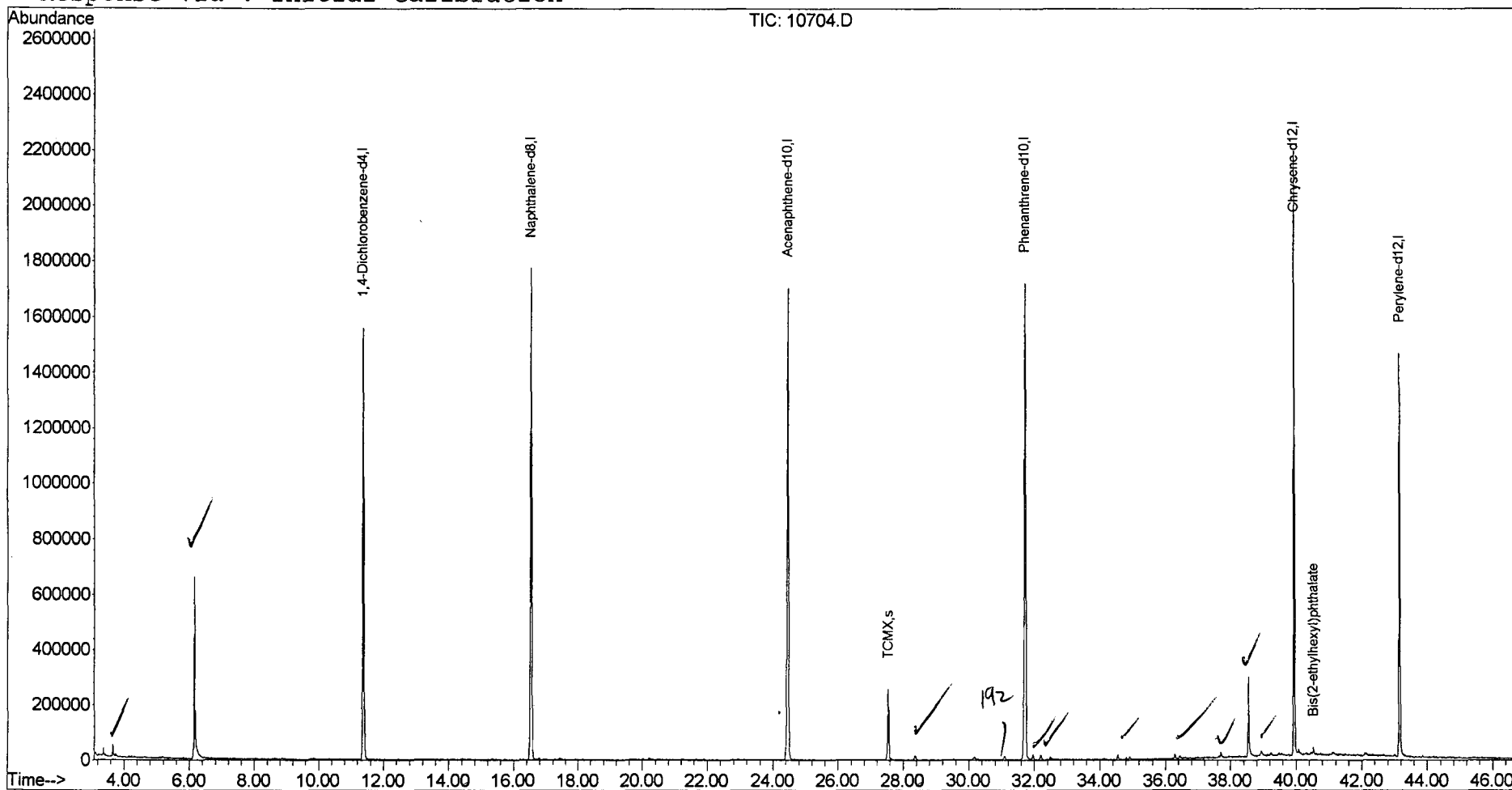
Quantitation Report (QT Reviewed)

Data File : C:\MSDCHEM\1\DATA\020513\10704.D  
 Acq On : 14 May 2002 2:23 am  
 Sample : sample  
 Misc :  
 MS Integration Params: rteint.e  
 Quant Time: May 15 9:00 2002

Vial: 15  
 Operator: Nefliu  
 Inst : Instrumen  
 Multiplr: 1.00

Quant Results File: SCAN020507.RES

Method : C:\MSDCHEM\1\METHODS\SCAN020507.M (RTE Integrator)  
 Title : Method 508 GC/MS Scan  
 Last Update : Thu May 09 09:21:01 2002  
 Response via : Initial Calibration



Data File : C:\MSDCHEM\1\DATA\020513\10704.D

Acq On : 14 May 2002 2:23 am

Sample : sample

Misc :

MS Integration Params: rteint.e

Vial: 15

Operator: Nefliu

Inst : Instrumen

Multiplr: 1.00

Method : C:\MSDCHEM\1\METHODS\SCAN020507.M (RTE Integrator)

Title : Method 508 GC/MS Scan

Smoothing : ON

Sampling : 1

Start Thrs: 0.2

Stop Thrs : 0

Filtering: 5

Min Area: 0.7 % of largest Peak

Max Peaks: 100

Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >

Peak separation: 5

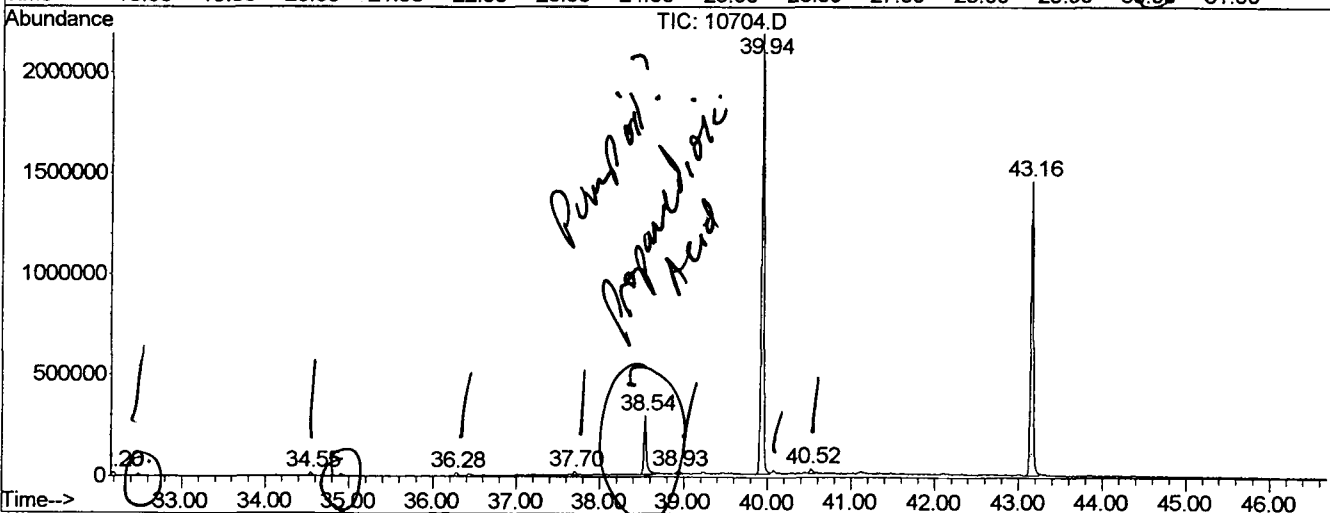
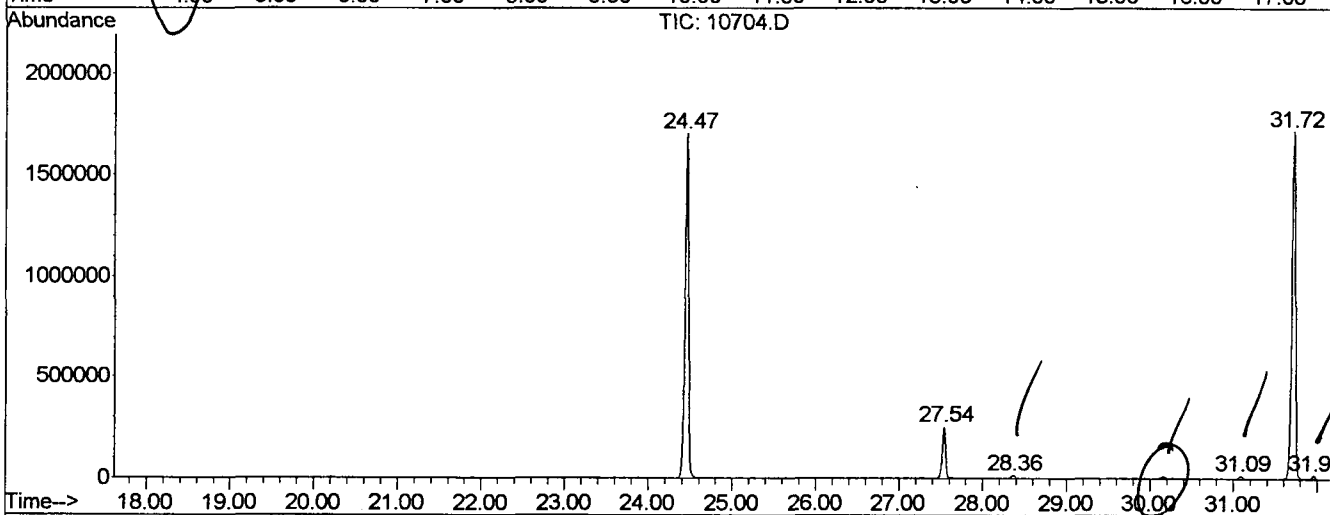
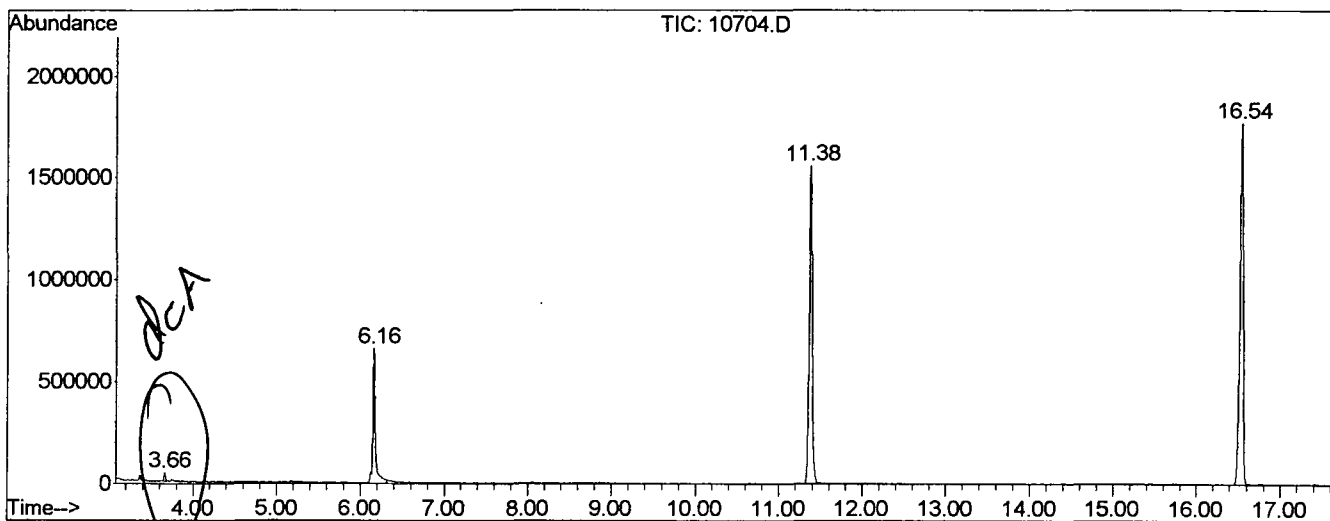
Signal : TIC

| 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         | 3.661       | 95            | 99          | 109          | rVB2     | 42472          | 54737         | 1.02%           | 0.183%        |
| 2         | 6.158       | 511           | 524         | 551          | rBV      | 659616         | 1371282       | 25.60%          | 4.591%        |
| 3         | 11.376      | 1400          | 1412        | 1436         | rBV      | 1556148        | 3888192       | 72.59%          | 13.019%       |
| 4         | 16.540      | 2273          | 2291        | 2308         | rBV      | 1770658        | 4849100       | 90.53%          | 16.236%       |
| 5         | 24.466      | 3617          | 3640        | 3657         | rBV2     | 1703080        | 5100813       | 95.23%          | 17.079%       |
| 6         | 27.539      | 4146          | 4163        | 4175         | rBV2     | 254524         | 699480        | 13.06%          | 2.342%        |
| 7         | 28.356      | 4291          | 4302        | 4310         | rBV4     | 16832          | 54873         | 1.02%           | 0.184%        |
| 8         | 31.088      | 4757          | 4767        | 4777         | rBV4     | 14894          | 46709         | 0.87%           | 0.156%        |
| 9         | 31.723      | 4855          | 4875        | 4900         | rVB3     | 1715415        | 5356371       | 100.00%         | 17.935%       |
| 10        | 31.958      | 4906          | 4915        | 4924         | rVB4     | 17904          | 48682         | 0.91%           | 0.163%        |
| 11        | 32.199      | 4942          | 4956        | 4966         | rBV2     | 17983          | 53976         | 1.01%           | 0.181%        |
| 12        | 34.549      | 5343          | 5356        | 5363         | rBV2     | 17759          | 40107         | 0.75%           | 0.134%        |
| 13        | 36.276      | 5643          | 5650        | 5660         | rBV4     | 16916          | 43002         | 0.80%           | 0.144%        |
| 14        | 37.698      | 5886          | 5892        | 5902         | rBV7     | 17169          | 42649         | 0.80%           | 0.143%        |
| 15        | 38.544      | 6027          | 6036        | 6057         | rBV2     | 289838         | 647211        | 12.08%          | 2.167%        |
| 16        | 38.932      | 6096          | 6102        | 6115         | rBV8     | 14594          | 42673         | 0.80%           | 0.143%        |
| 17        | 39.937      | 6262          | 6273        | 6288         | rBV2     | 2175475        | 4355358       | 81.31%          | 14.583%       |
| 18        | 40.518      | 6367          | 6372        | 6380         | rVB3     | 25611          | 54107         | 1.01%           | 0.181%        |
| 19        | 43.157      | 6809          | 6821        | 6836         | rBV2     | 1458042        | 3116626       | 58.19%          | 10.435%       |

Sum of corrected areas: 29865948

LSC Report - Integrated Chromatogram

File : C:\MSDCHEM\1\DATA\020513\10704.D  
 Operator : Nefliu  
 Acquired : 14 May 2002 2:23 am using AcqMethod 508SCAN1  
 Instrument : Instrumen  
 Sample Name: sample  
 Misc Info :  
 Vial Number: 15  
 Quant File :SCAN020507.RES (RTE Integrator)



Data File : C:\MSDCHEM\1\DATA\020513\10704.D  
 Acq On : 14 May 2002 2:23 am  
 Sample : sample  
 Misc :  
 MS Integration Params: rteint.e

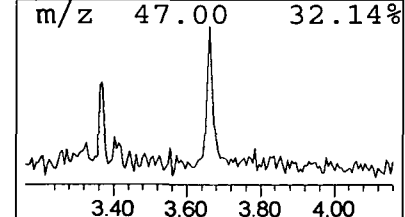
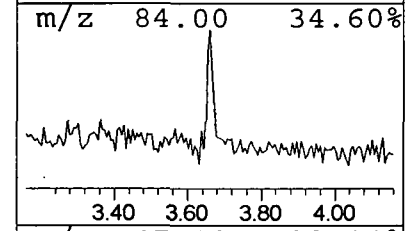
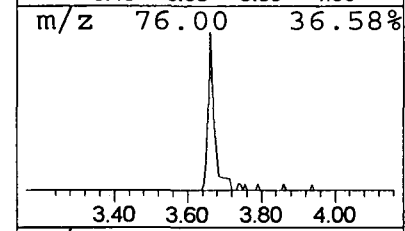
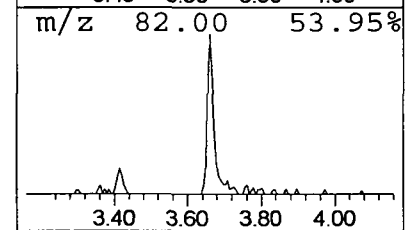
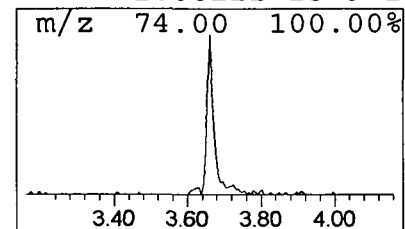
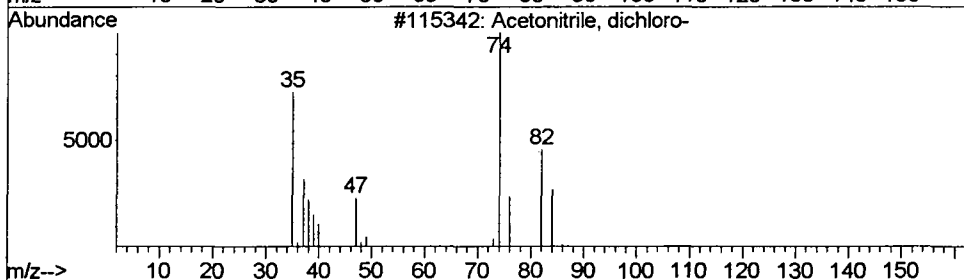
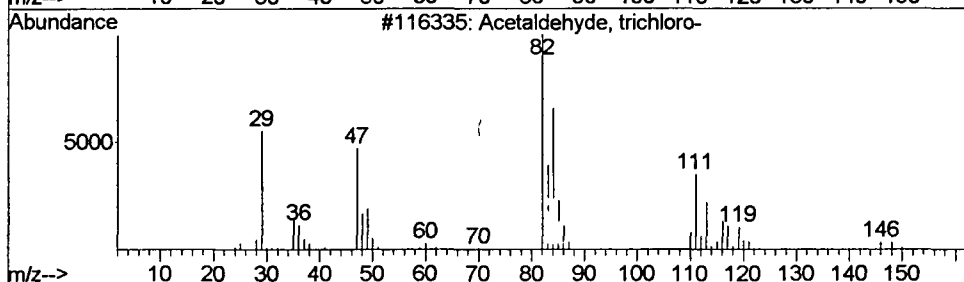
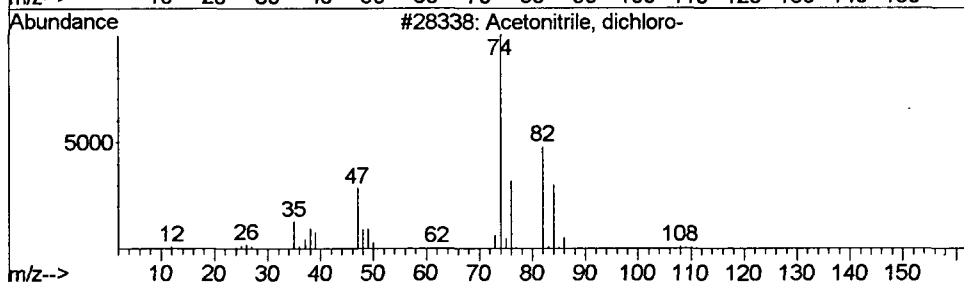
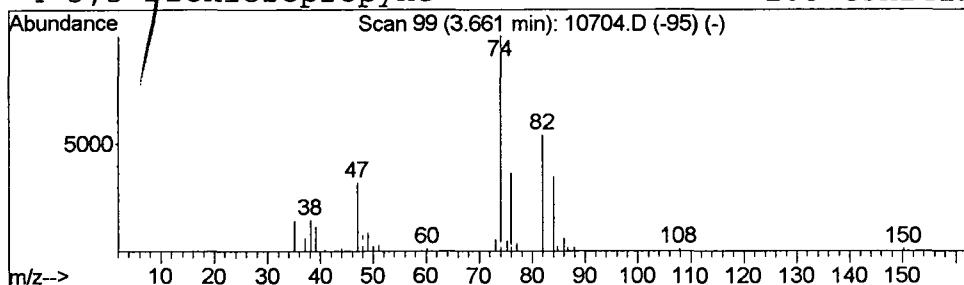
Vial: 15  
 Operator: Nefliu  
 Inst : Instrumen  
 Multiplr: 1.00

Quant Method : C:\MSDCHEM\1\METHODS\SCAN020507.M (RTE Integrator)  
 Title : Method 508 GC/MS Scan  
 Library : C:\DATABASE\nist98.1

\*\*\*\*\*  
 Peak Number 1 Acetonitrile, dichloro- Concentration Rank 3

| R.T. | EstConc   | Area  | Relative to ISTD       | R.T.  |
|------|-----------|-------|------------------------|-------|
| 3.66 | 0.56 ug/l | 54737 | 1,4-Dichlorobenzene-d4 | 11.38 |

| Hit# of 5 | Tentative ID             | MW  | MolForm | CAS#         | Qual |
|-----------|--------------------------|-----|---------|--------------|------|
| 1         | Acetonitrile, dichloro-  | 109 | C2HCl2N | 003018-12-0  | 91   |
| 2         | Acetaldehyde, trichloro- | 146 | C2HCl3O | 000075-87-6  | 38   |
| 3         | Acetonitrile, dichloro-  | 109 | C2HCl2N | 003018-12-0  | 37   |
| 4         | 3,3-Dichloropropyne      | 108 | C3H2Cl2 | 1000222-23-9 | 28   |



Data File : C:\MSDCHEM\1\DATA\020513\10704.D  
 Acq On : 14 May 2002 2:23 am  
 Sample : sample  
 Misc :  
 MS Integration Params: rteint.e

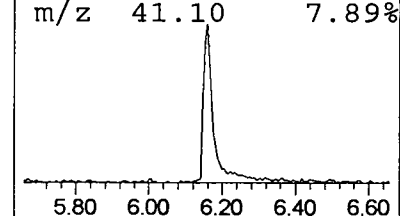
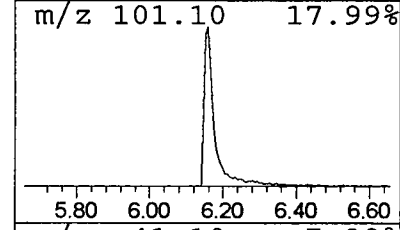
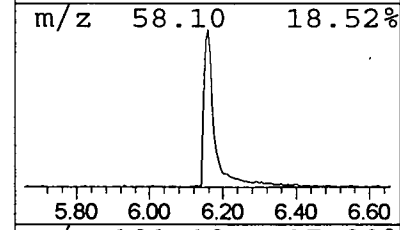
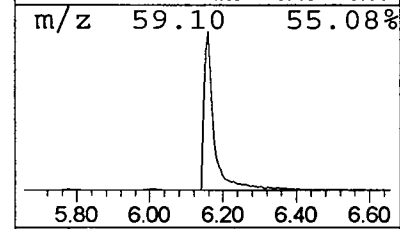
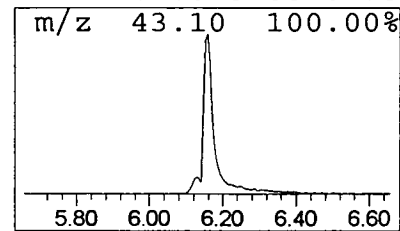
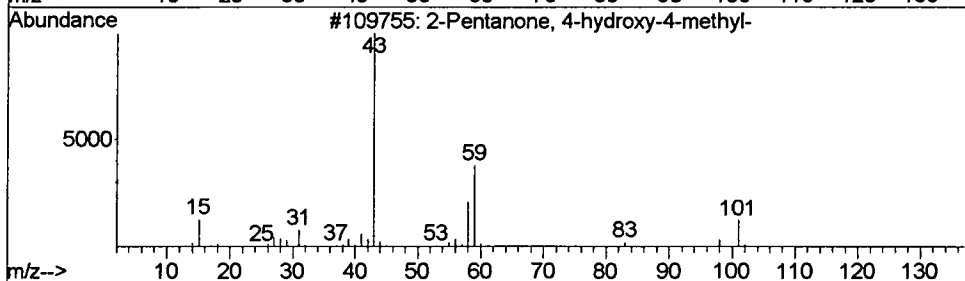
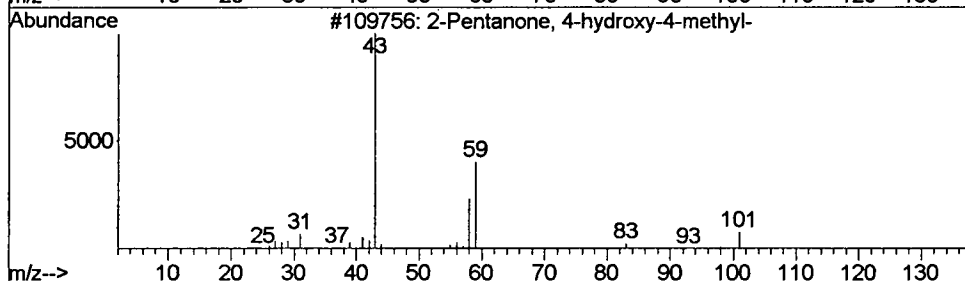
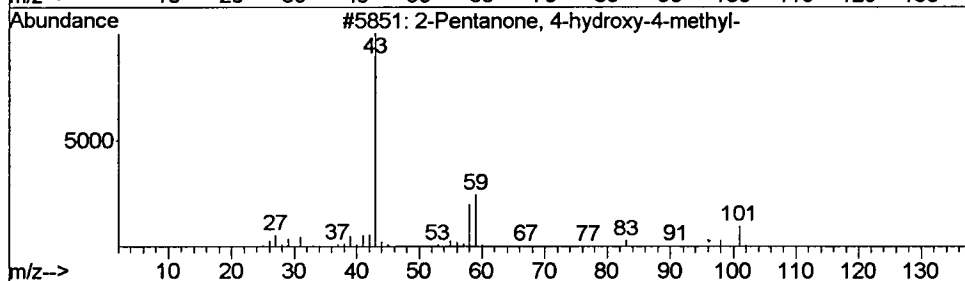
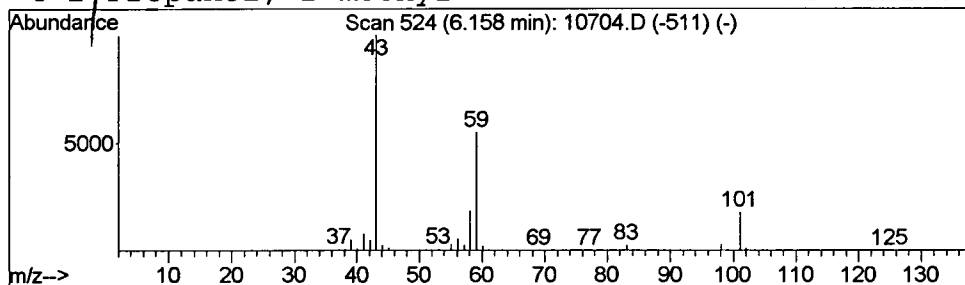
Vial: 15  
 Operator: Nefliu  
 Inst : Instrumen  
 Multiplr: 1.00

Quant Method : C:\MSDCHEM\1\METHODS\SCAN020507.M (RTE Integrator)  
 Title : Method 508 GC/MS Scan  
 Library : C:\DATABASE\nist98.l

\*\*\*\*\*  
 Peak Number 2 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 1

| R.T. | EstConc    | Area    | Relative to ISTD       | R.T.  |
|------|------------|---------|------------------------|-------|
| 6.16 | 14.11 ug/l | 1371280 | 1,4-Dichlorobenzene-d4 | 11.38 |

| Hit# | of | 5 | Tentative ID                     | MW  | MolForm | CAS#        | Qual |
|------|----|---|----------------------------------|-----|---------|-------------|------|
| ✓ 1  |    |   | 2-Pentanone, 4-hydroxy-4-methyl- | 116 | C6H12O2 | 000123-42-2 | 50   |
| 2    |    |   | 2-Pentanone, 4-hydroxy-4-methyl- | 116 | C6H12O2 | 000123-42-2 | 45   |
| 3    |    |   | 2-Pentanone, 4-hydroxy-4-methyl- | 116 | C6H12O2 | 000123-42-2 | 45   |
| 4    |    |   | 2-Propanol, 2-methyl-            | 74  | C4H10O  | 000075-65-0 | 25   |



Data File : C:\MSDCHEM\1\DATA\020513\10704.D  
Acq On : 14 May 2002 2:23 am  
Sample : sample  
Misc :  
MS Integration Params: rteint.e

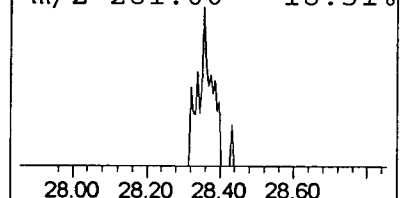
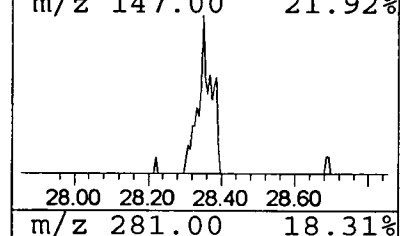
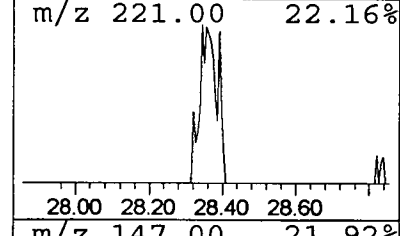
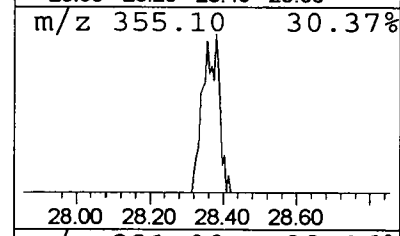
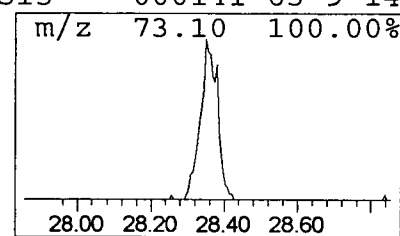
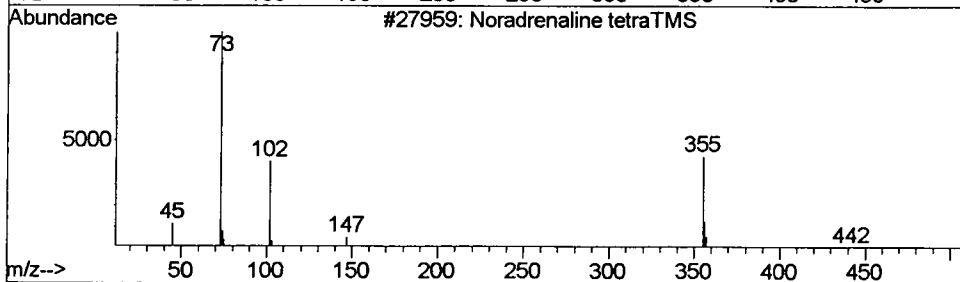
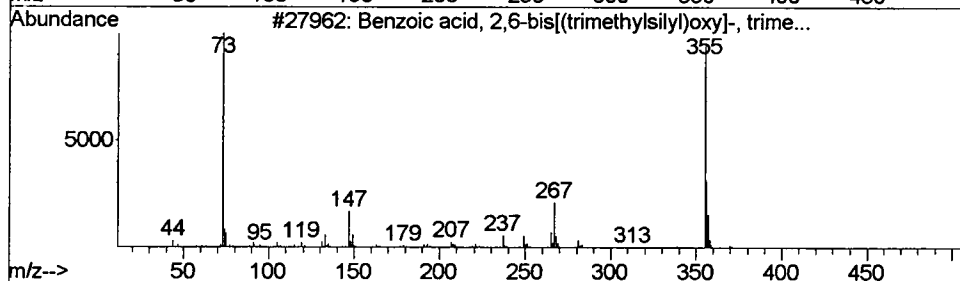
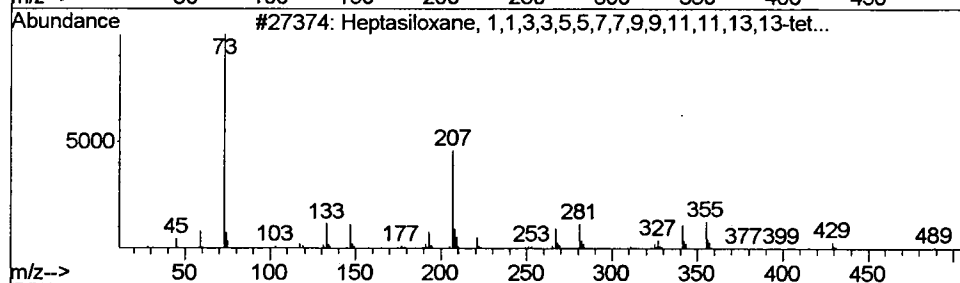
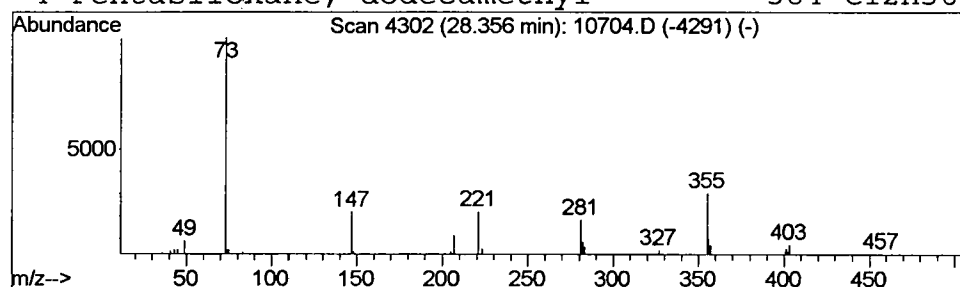
Vial: 15  
Operator: Nefliu  
Inst : Instrumen  
Multiplr: 1.00

Quant Method : C:\MSDCHEM\1\METHODS\SCAN020507.M (RTE Integrator)  
Title : Method 508 GC/MS Scan  
Library : C:\DATABASE\nist98.1

\*\*\*\*\*  
Peak Number 3 Heptasiloxane, 1,1,3,3,5,5,... Concentration Rank 4

| R.T.  | EstConc   | Area  | Relative to ISTD | R.T.  |
|-------|-----------|-------|------------------|-------|
| 28.36 | 0.41 ug/l | 54873 | Phenanthrene-d10 | 31.72 |

| Hit# of | 5                                   | Tentative ID | MW           | MolForm     | CAS# | Qual |
|---------|-------------------------------------|--------------|--------------|-------------|------|------|
| 1       | Heptasiloxane, 1,1,3,3,5,5,7,7,9... | 504          | C14H44O6Si7  | 019095-23-9 | 23   |      |
| 2       | Benzoic acid, 2,6-bis[(trimethyl... | 370          | C16H30O4Si3  | 003782-85-2 | 23   |      |
| 3       | Noradrenaline tetraTMS              | 457          | C20H43NO3Si4 | 068595-65-3 | 17   |      |
| 4       | Pentasiloxane, dodecamethyl-        | 384          | C12H36O4Si5  | 000141-63-9 | 14   |      |



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 Acq On : 14 May 2002 2:23 am  
 Sample : sample  
 Misc :  
 MS Integration Params: rteint.e

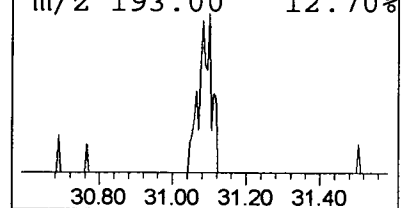
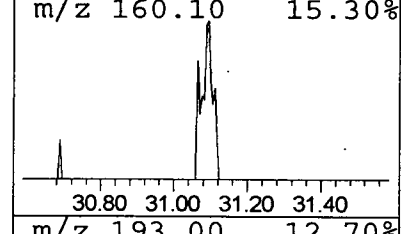
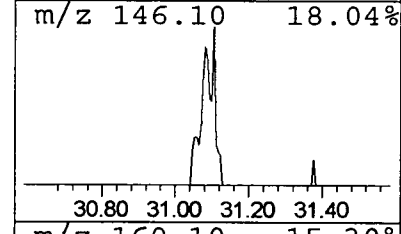
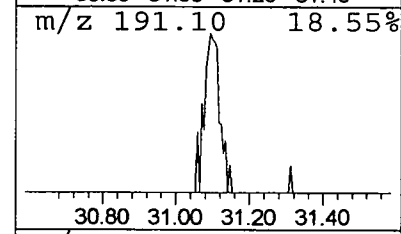
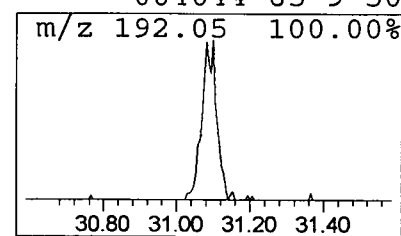
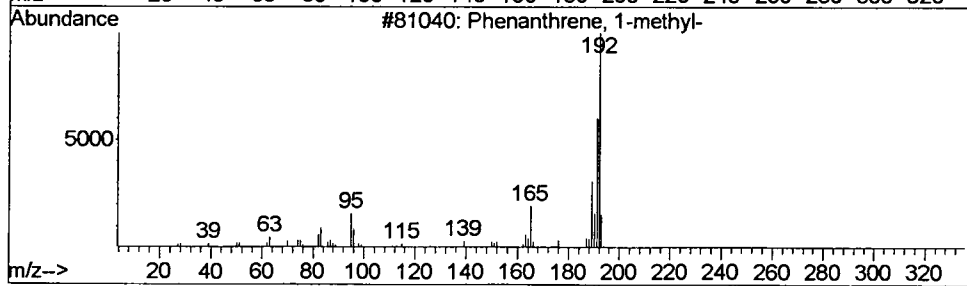
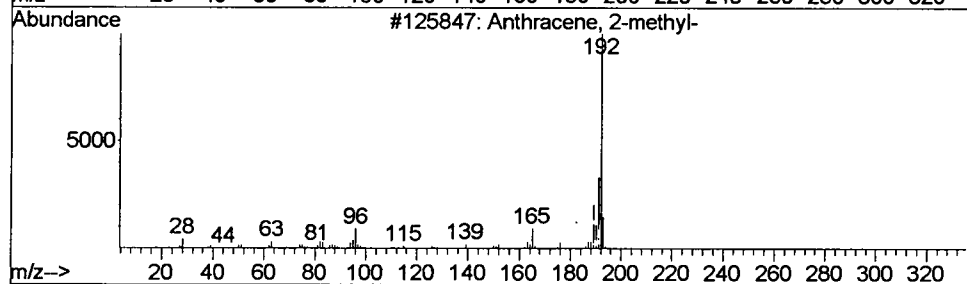
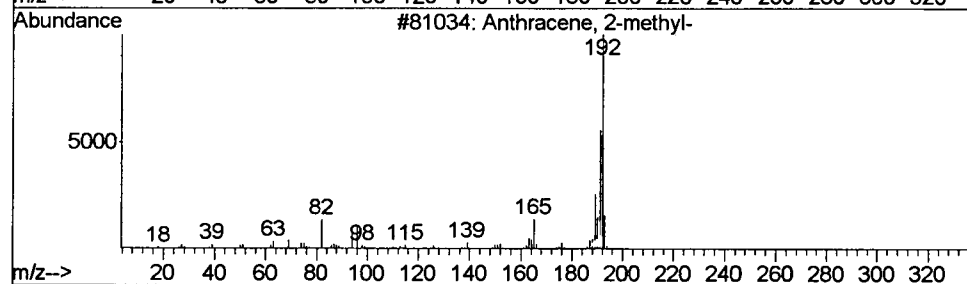
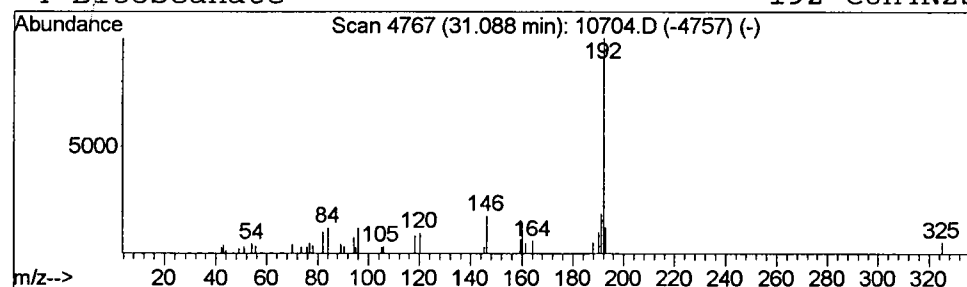
Vial: 15  
 Operator: Nefliu  
 Inst : Instrumen  
 Multiplr: 1.00

Quant Method : C:\MSDCHEM\1\METHODS\SCAN020507.M (RTE Integrator)  
 Title : Method 508 GC/MS Scan  
 Library : C:\DATABASE\nist98.1

\*\*\*\*\*  
 Peak Number 4 Anthracene, 2-methyl- Concentration Rank 10

| R.T.  | EstConc   | Area  | Relative to ISTD | R.T.  |
|-------|-----------|-------|------------------|-------|
| 31.09 | 0.35 ug/l | 46709 | Phenanthrene-d10 | 31.72 |

| Hit# | of | Tentative ID            | MW  | MolForm  | CAS#        | Qual |
|------|----|-------------------------|-----|----------|-------------|------|
| 1    | 5  | Anthracene, 2-methyl-   | 192 | C15H12   | 000613-12-7 | 58   |
| 2    |    | Anthracene, 2-methyl-   | 192 | C15H12   | 000613-12-7 | 53   |
| 3    |    | Phenanthrene, 1-methyl- | 192 | C15H12   | 000832-69-9 | 50   |
| 4    |    | Bitoscanate             | 192 | C8H4N2S2 | 004044-65-9 | 50   |



Data File : C:\MSDCHEM\1\DATA\020513\10704.D

Acq On : 14 May 2002 2:23 am

Sample : sample

Misc :

MS Integration Params: rteint.e

Vial: 15

Operator: Nefliu

Inst : Instrumen

Multiplr: 1.00

Quant Method : C:\MSDCHEM\1\METHODS\SCAN020507.M (RTE Integrator)

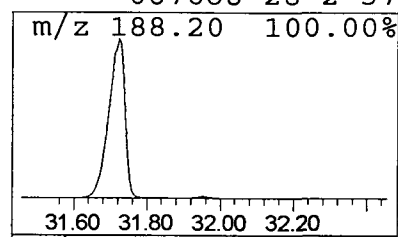
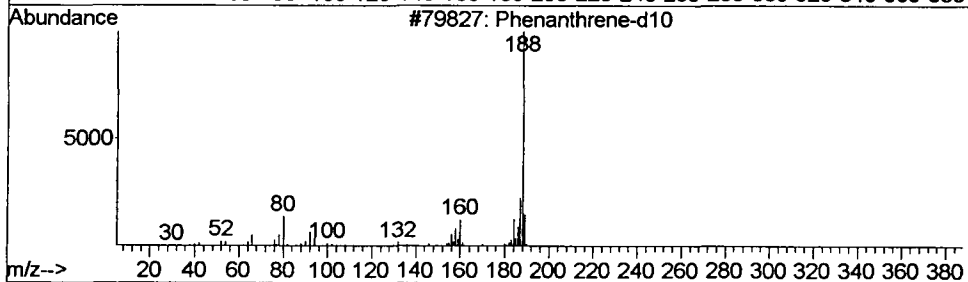
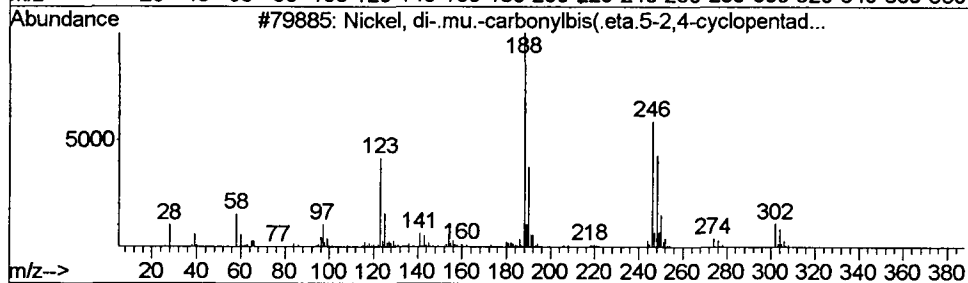
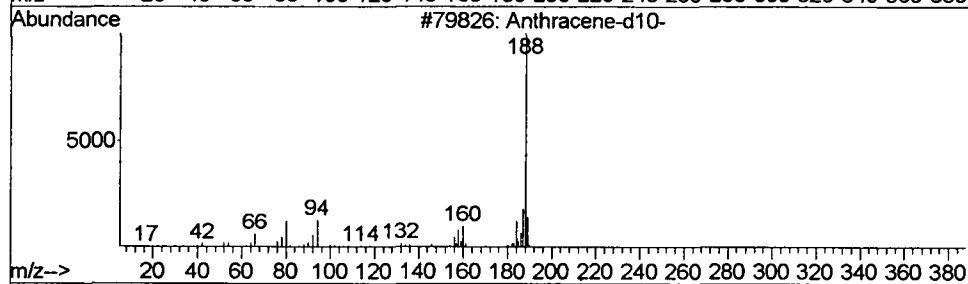
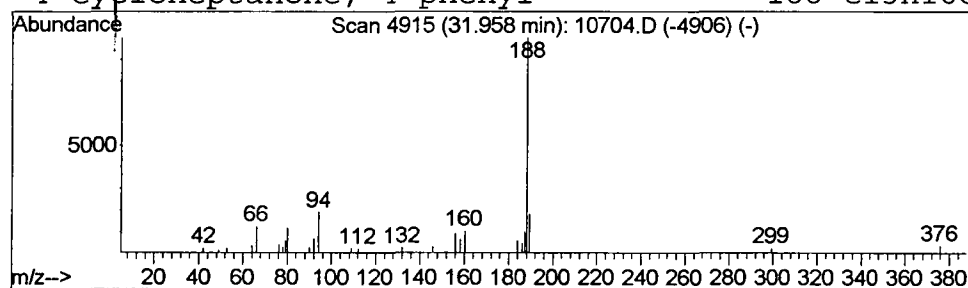
Title : Method 508 GC/MS Scan

Library : C:\DATABASE\mist98.1

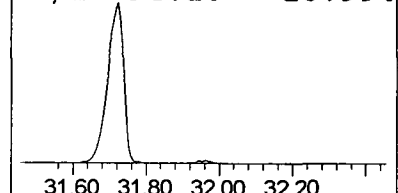
\*\*\*\*\*  
Peak Number 5 Anthracene-d10- Concentration Rank 9

| R.T.  | EstConc   | Area  | Relative to ISTD | R.T.  |
|-------|-----------|-------|------------------|-------|
| 31.96 | 0.36 ug/l | 48682 | Phenanthrene-d10 | 31.72 |

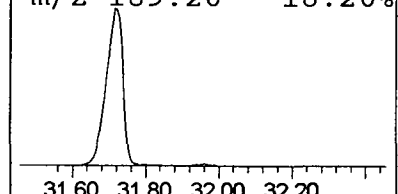
| Hit# | of | 5 | Tentative ID                        | MW  | MolForm     | CAS#        | Qual |
|------|----|---|-------------------------------------|-----|-------------|-------------|------|
| ✓ 1  |    |   | Anthracene-d10-                     | 188 | C14D10      | 001719-06-8 | 72   |
| 2    |    |   | Nickel, di-.mu.-carbonylbis(.eta... | 302 | C12H10Ni2O2 | 012170-92-2 | 43   |
| 3    |    |   | Phenanthrene-d10                    | 188 | C14D10      | 001517-22-2 | 39   |
| 4    |    |   | Cycloheptanone, 4-phenyl-           | 188 | C13H16O     | 067688-28-2 | 37   |



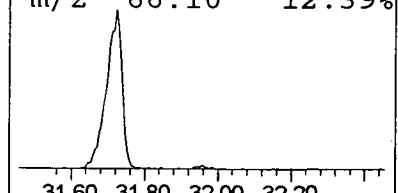
m/z 94.10 18.99%



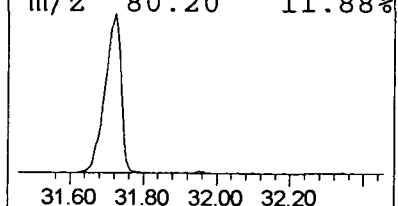
m/z 189.20 18.20%



m/z 66.10 12.39%



m/z 80.20 11.88%



Data File : C:\MSDCHEM\1\DATA\020513\10704.D

Acq On : 14 May 2002 2:23 am

Sample : sample

Misc :

MS Integration Params: rteint.e

Vial: 15

Operator: Nefliu

Inst : Instrumen

Multiplr: 1.00

Quant Method : C:\MSDCHEM\1\METHODS\SCAN020507.M (RTE Integrator)

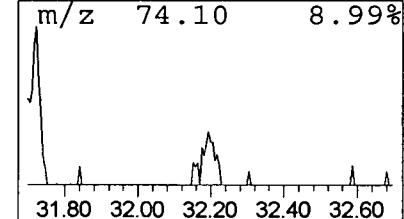
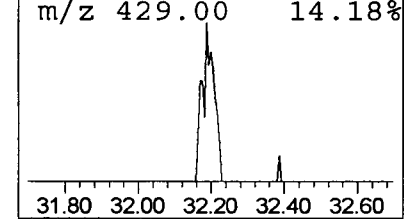
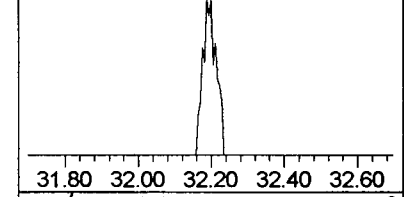
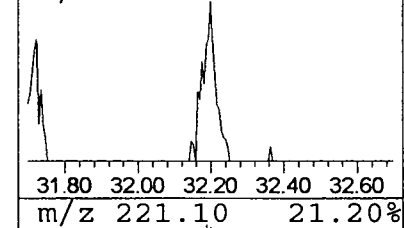
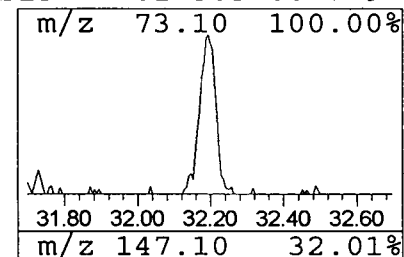
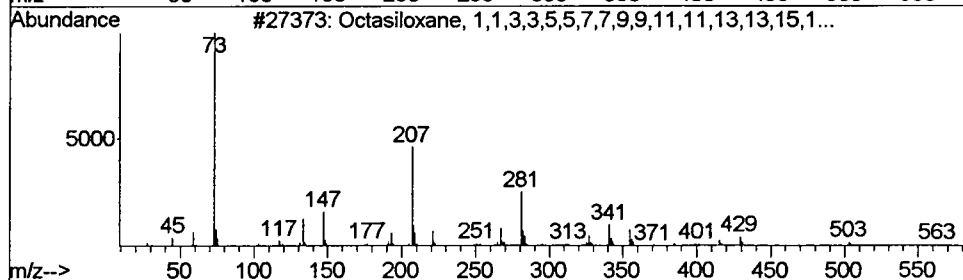
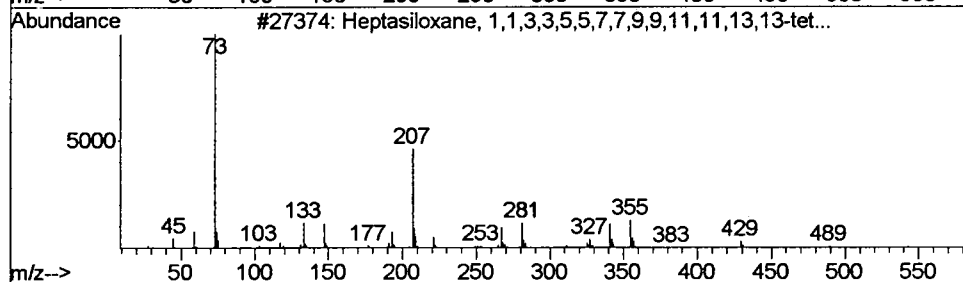
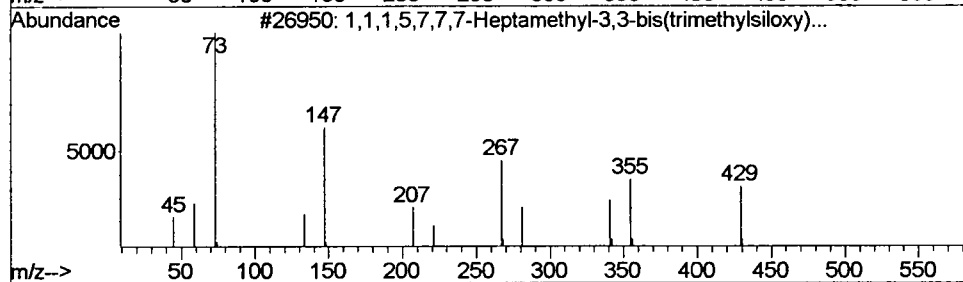
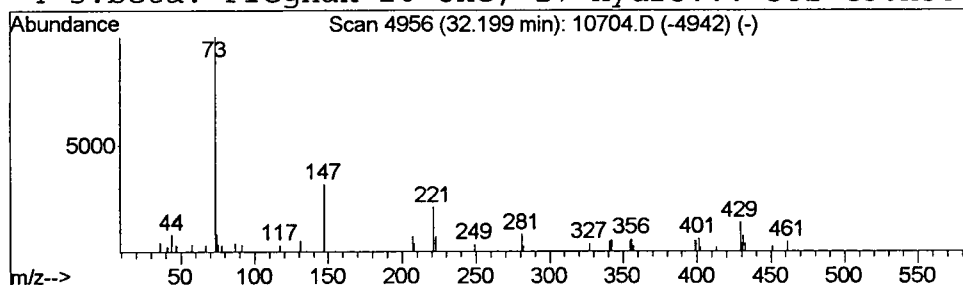
Title : Method 508 GC/MS Scan

Library : C:\DATABASE\nist98.l

\*\*\*\*\*  
Peak Number 6 1,1,1,5,7,7,7-Heptamethyl-3... Concentration Rank 5

| R.T.  | EstConc   | Area  | Relative to ISTD | R.T.  |
|-------|-----------|-------|------------------|-------|
| 32.20 | 0.40 ug/l | 53976 | Phenanthrene-d10 | 31.72 |

| Hit# | of                                  | 5   | Tentative ID | MW          | MolForm | CAS# | Qual |
|------|-------------------------------------|-----|--------------|-------------|---------|------|------|
| 1    | 1,1,1,5,7,7,7-Heptamethyl-3,3-bi... | 444 | C13H40O5Si6  | 038147-00-1 | 32      |      |      |
| 2    | Heptasiloxane, 1,1,3,3,5,5,7,7,9... | 504 | C14H44O6Si7  | 019095-23-9 | 10      |      |      |
| 3    | Octasiloxane, 1,1,3,3,5,5,7,7,9...  | 578 | C16H50O7Si8  | 019095-24-0 | 9       |      |      |
| 4    | 5.beta.-Pregnan-20-one, 17-hydro... | 582 | C30H58O5Si3  | 017563-00-7 | 9       |      |      |



Data File : C:\MSDCHEM\1\DATA\020513\10704.D  
 Acq On : 14 May 2002 2:23 am  
 Sample : sample  
 Misc :  
 MS Integration Params: rteint.e

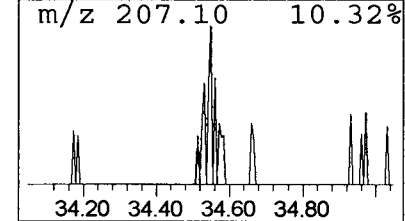
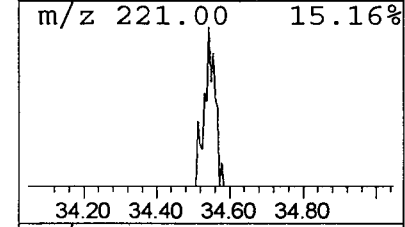
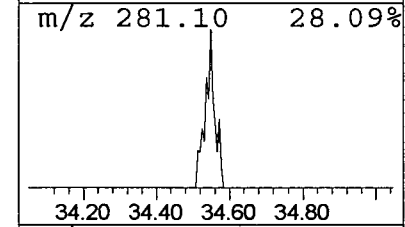
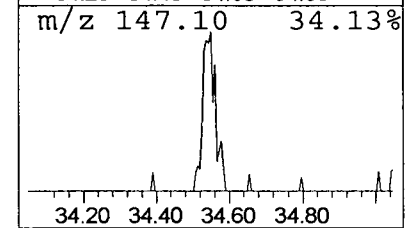
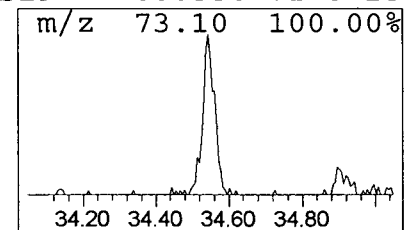
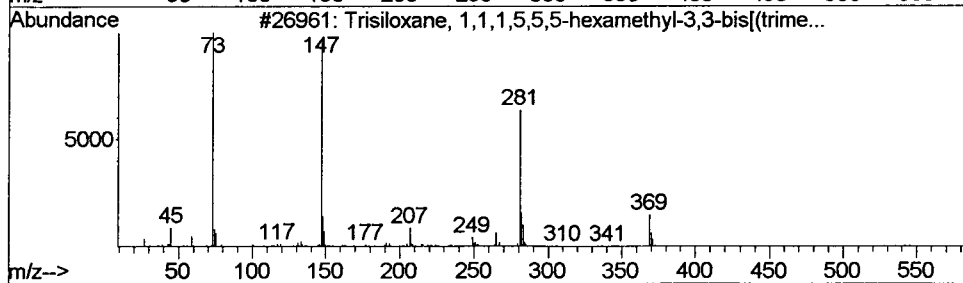
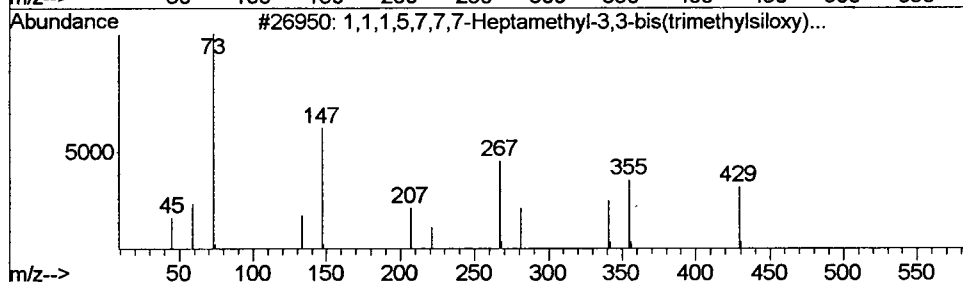
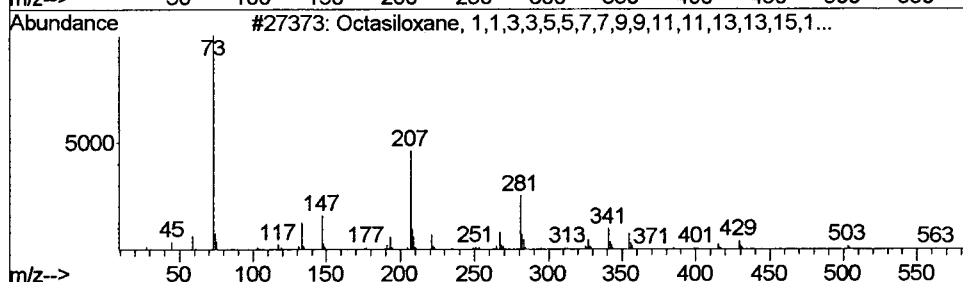
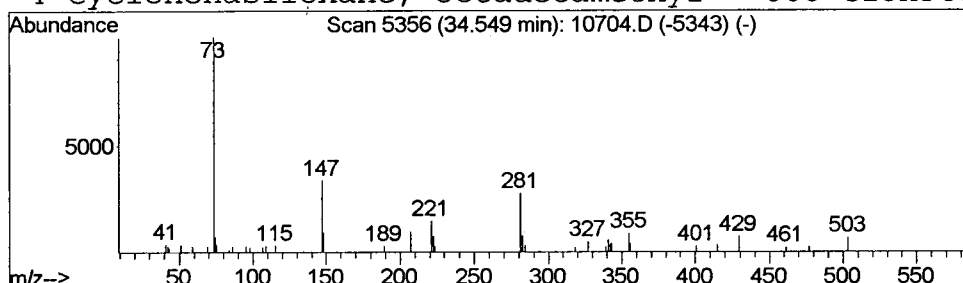
Vial: 15  
 Operator: Nefliu  
 Inst : Instrumen  
 Multiplr: 1.00

Quant Method : C:\MSDCHEM\1\METHODS\SCAN020507.M (RTE Integrator)  
 Title : Method 508 GC/MS Scan  
 Library : C:\DATABASE\nist98.1

\*\*\*\*\*  
 Peak Number 7 Octasiloxane, 1,1,3,3,5,5,7... Concentration Rank 11

| R.T.  | EstConc   | Area  | Relative to ISTD | R.T.  |
|-------|-----------|-------|------------------|-------|
| 34.55 | 0.30 ug/l | 40107 | Phenanthrene-d10 | 31.72 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm     | CAS#        | Qual |
|------|----|---|-------------------------------------|-----|-------------|-------------|------|
| 1    |    |   | Octasiloxane, 1,1,3,3,5,5,7,7,9...  | 578 | C16H50O7Si8 | 019095-24-0 | 45   |
| 2    |    |   | 1,1,1,5,7,7,7-Heptamethyl-3,3-bi... | 444 | C13H40O5Si6 | 038147-00-1 | 37   |
| 3    |    |   | Trisiloxane, 1,1,1,5,5,5-hexamet... | 384 | C12H36O4Si5 | 003555-47-3 | 33   |
| 4    |    |   | Cyclononasiloxane, octadecamethyl-  | 666 | C18H54O9Si9 | 000556-71-8 | 25   |



Data File : C:\MSDCHEM\1\DATA\020513\10704.D  
 Acq On : 14 May 2002 2:23 am  
 Sample : sample  
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 MS Integration Params: rteint.e

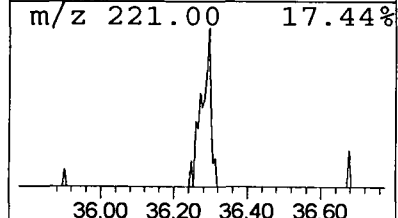
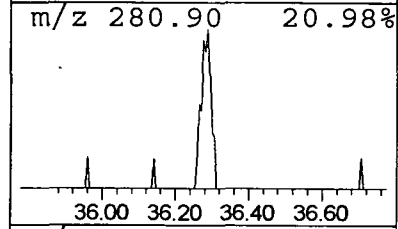
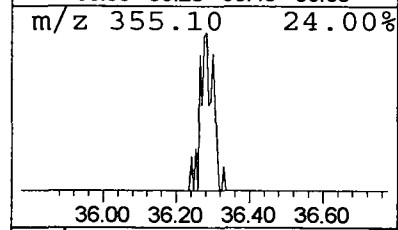
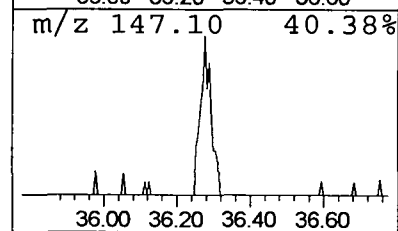
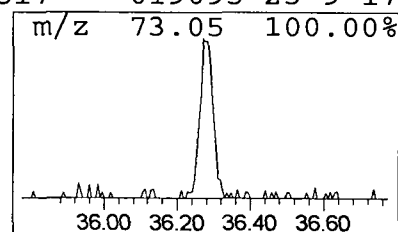
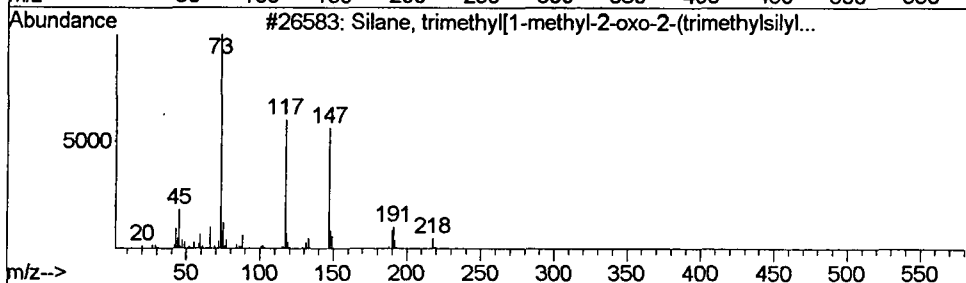
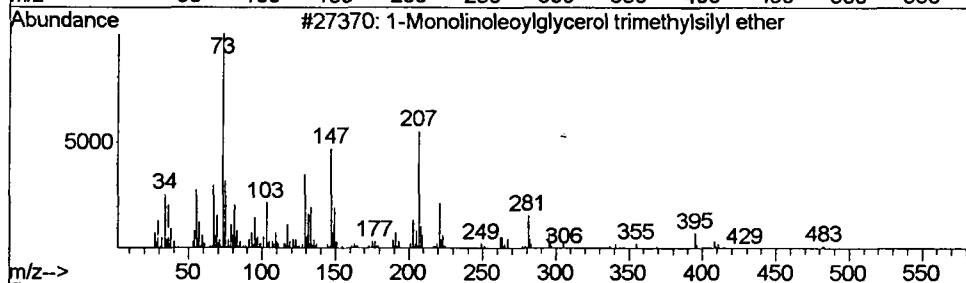
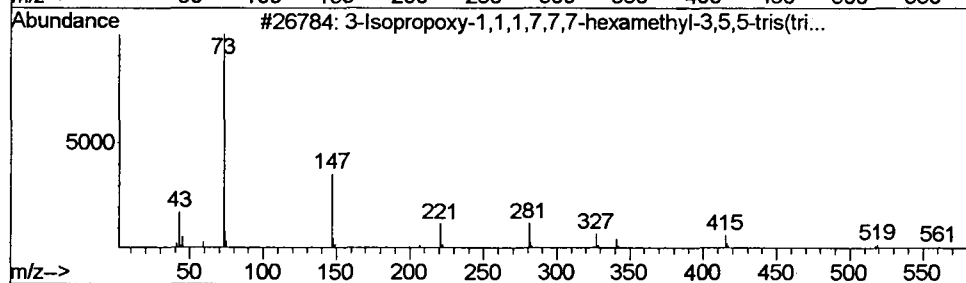
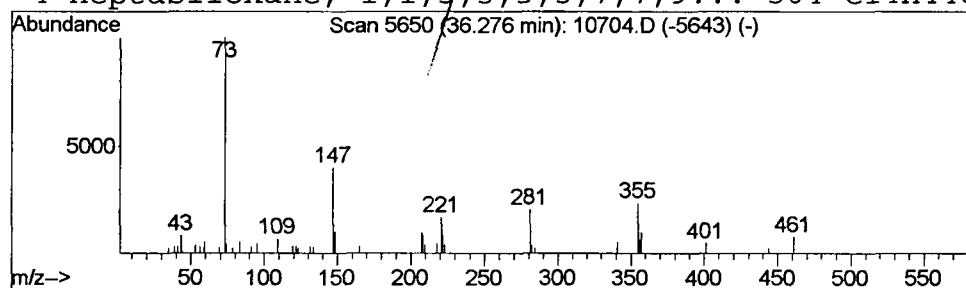
Vial: 15  
 Operator: Nefliu  
 Inst : Instrumen  
 Multiplr: 1.00

Quant Method : C:\MSDCHEM\1\METHODS\SCAN020507.M (RTE Integrator)  
 Title : Method 508 GC/MS Scan  
 Library : C:\DATABASE\nist98.1

\*\*\*\*\*  
 Peak Number 8 3-Isopropoxy-1,1,1,7,7,7-he... Concentration Rank 6

| R.T.  | EstConc   | Area  | Relative to ISTD | R.T.  |
|-------|-----------|-------|------------------|-------|
| 36.28 | 0.39 ug/l | 43002 | Chrysene-d12     | 39.94 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm     | CAS#        | Qual |
|------|----|---|-------------------------------------|-----|-------------|-------------|------|
| 1    |    |   | 3-Isopropoxy-1,1,1,7,7,7-hexamet... | 576 | C18H52O7Si7 | 071579-69-6 | 38   |
| 2    |    |   | 1-Monolinoleoylglycerol trimethy... | 498 | C27H54O4Si2 | 054284-45-6 | 28   |
| 3    |    |   | Silane, trimethyl[1-methyl-2-oxo... | 218 | C9H22O2Si2  | 055255-93-1 | 27   |
| 4    |    |   | Heptasiloxane, 1,1,3,3,5,5,7,7,9... | 504 | C14H44O6Si7 | 019095-23-9 | 17   |



Data File : C:\MSDCHEM\1\DATA\020513\10704.D

Acq On : 14 May 2002 2:23 am

Sample : sample

Misc :

MS Integration Params: rteint.e

Vial: 15

Operator: Nefliu

Inst : Instrumen

Multiplr: 1.00

Quant Method : C:\MSDCHEM\1\METHODS\SCAN020507.M (RTE Integrator)

Title : Method 508 GC/MS Scan

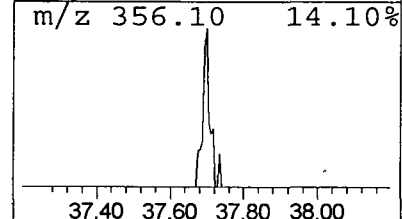
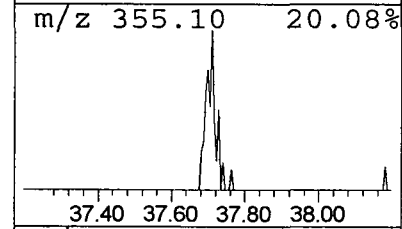
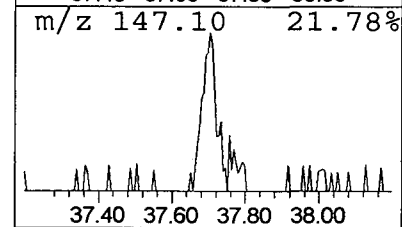
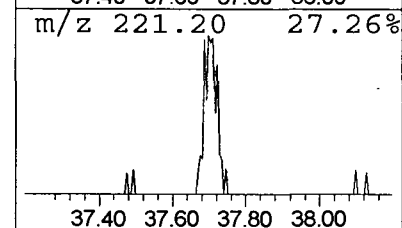
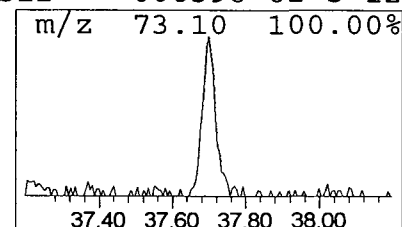
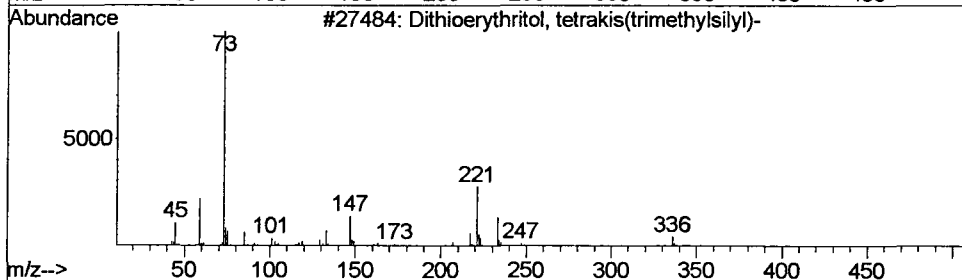
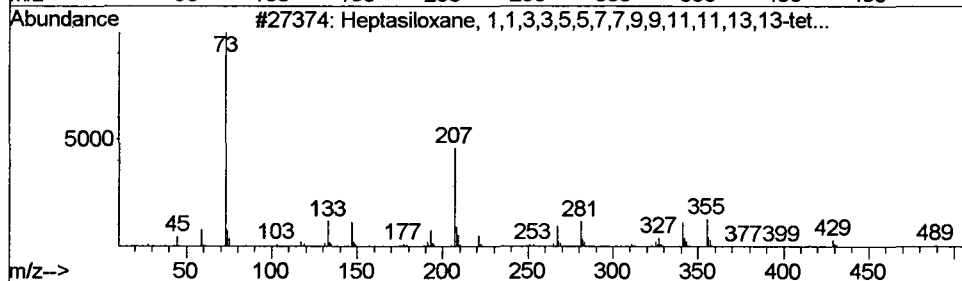
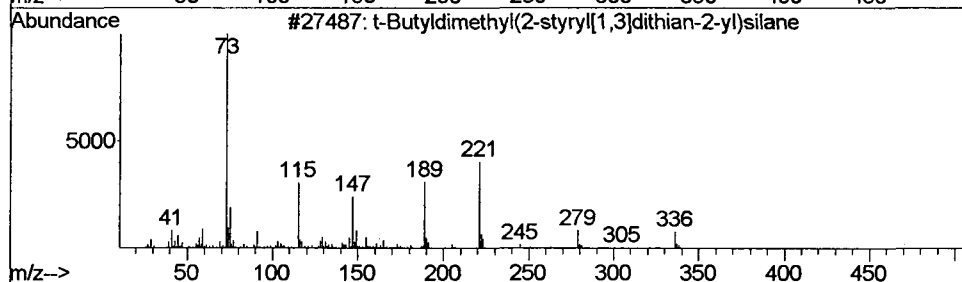
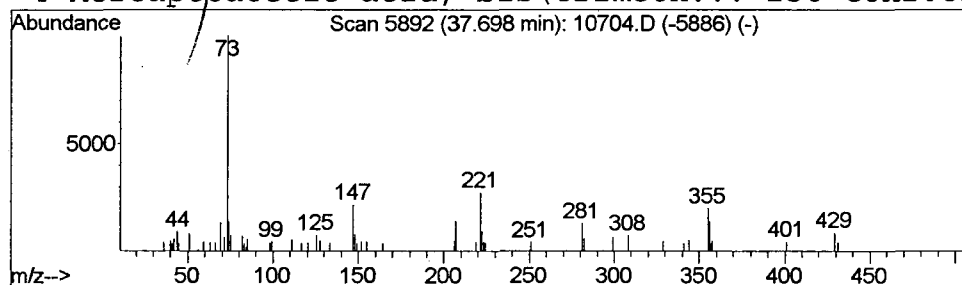
Library : C:\DATABASE\nist98.1

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Peak Number 9 t-Butyldimethyl(2-styryl[1,... Concentration Rank 8

| R.T.  | EstConc   | Area  | Relative to ISTD | R.T.  |
|-------|-----------|-------|------------------|-------|
| 37.70 | 0.39 ug/l | 42649 | Chrysene-d12     | 39.94 |

| Hit# | of | 5 | Tentative ID                          | MW  | MolForm.      | CAS#         | Qual |
|------|----|---|---------------------------------------|-----|---------------|--------------|------|
| 1    |    |   | t-Butyldimethyl(2-styryl[1,3]dit...   | 336 | C18H28S2Si    | 1000192-03-8 | 37   |
| 2    |    |   | Heptasiloxane, 1,1,3,3,5,5,7,7,9...   | 504 | C14H44O6Si7   | 019095-23-9  | 23   |
| 3    |    |   | Dithioerythritol, tetrakis(trimeth... | 442 | C16H42O2S2Si4 | 1000079-30-7 | 17   |
| 4    |    |   | Mercaptacetic acid, bis(trimeth...    | 236 | C8H20O2SSi2   | 006398-62-5  | 12   |



Data File : C:\MSDCHEM\1\DATA\020513\10704.D  
Acq On : 14 May 2002 2:23 am  
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Misc :  
MS Integration Params: rteint.e

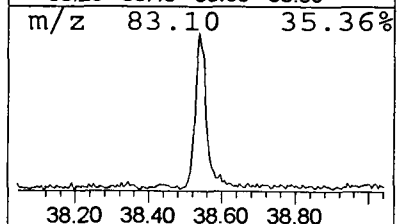
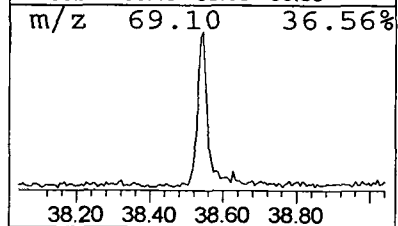
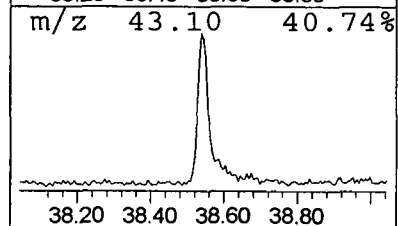
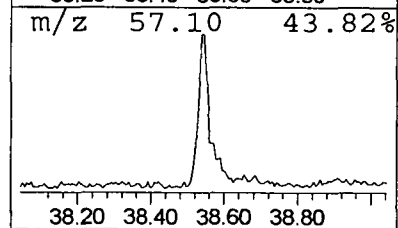
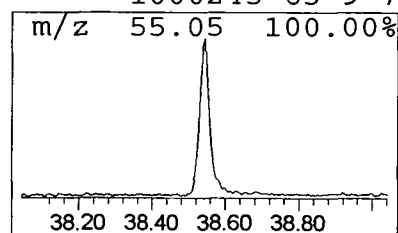
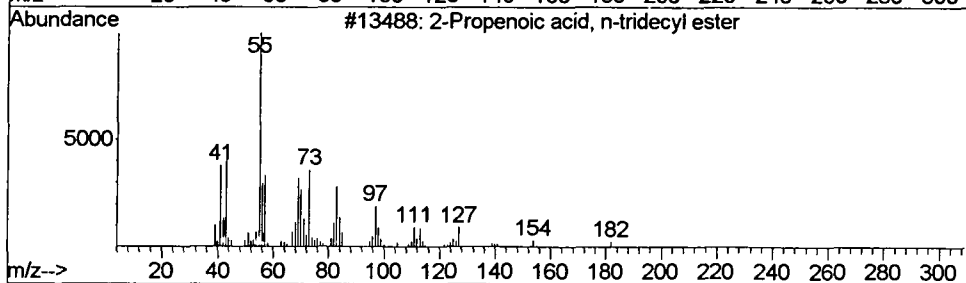
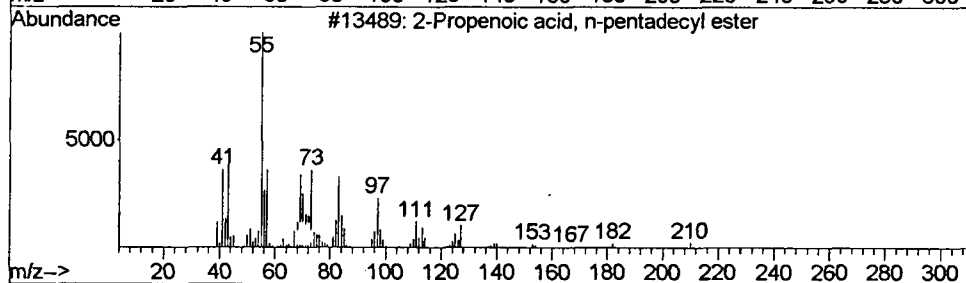
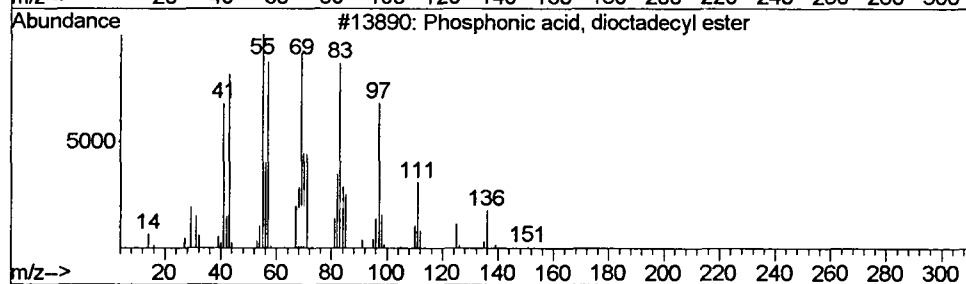
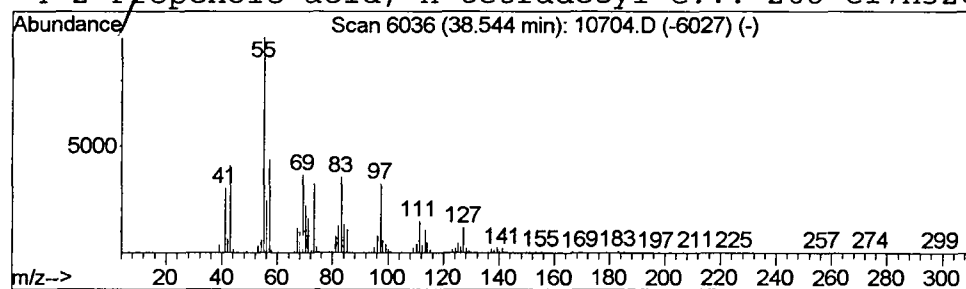
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Operator: Nefliu  
Inst : Instrumen  
Multiplr: 1.00

Quant Method : C:\MSDCHEM\1\METHODS\SCAN020507.M (RTE Integrator)  
Title : Method 508 GC/MS Scan  
Library : C:\DATABASE\nist98.1

\*\*\*\*\*  
Peak Number 10 Phosphonic acid, dioctadecy... Concentration Rank 2

| R.T.  | EstConc   | Area   | Relative to ISTD | R.T.  |
|-------|-----------|--------|------------------|-------|
| 38.54 | 5.94 ug/l | 647211 | Chrysene-d12     | 39.94 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm   | CAS#         | Qual |
|------|----|---|-------------------------------------|-----|-----------|--------------|------|
| 1    |    |   | Phosphonic acid, dioctadecyl ester  | 587 | C36H75O3P | 019047-85-9  | 91   |
| 2    |    |   | 2-Propenoic acid, n-pentadecyl e... | 282 | C18H34O2  | 1000245-64-0 | 90   |
| 3    |    |   | 2-Propenoic acid, n-tridecyl ester  | 254 | C16H30O2  | 1000245-63-8 | 87   |
| 4    |    |   | 2-Propenoic acid, n-tetradecyl e... | 268 | C17H32O2  | 1000245-63-9 | 72   |



Data File : C:\MSDCHEM\1\DATA\020513\10704.D  
 Acq On : 14 May 2002 2:23 am  
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 MS Integration Params: rteint.e

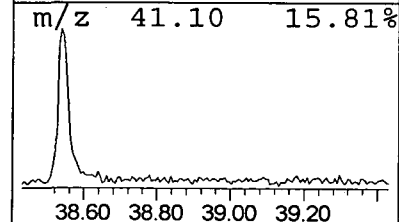
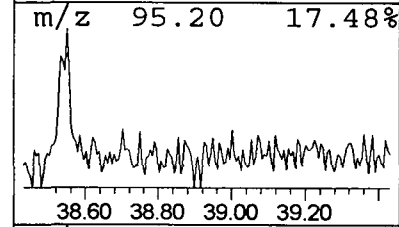
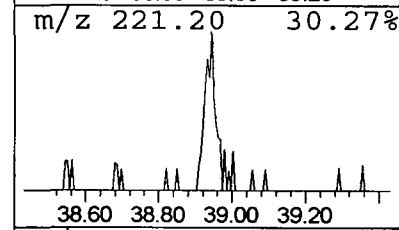
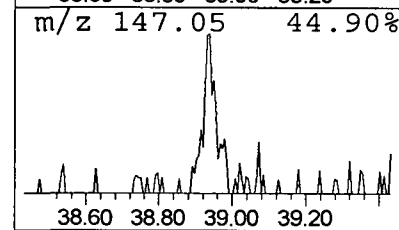
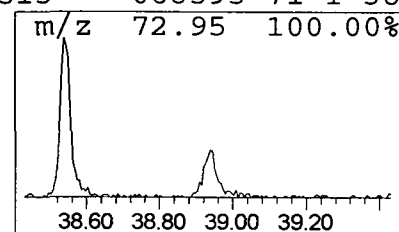
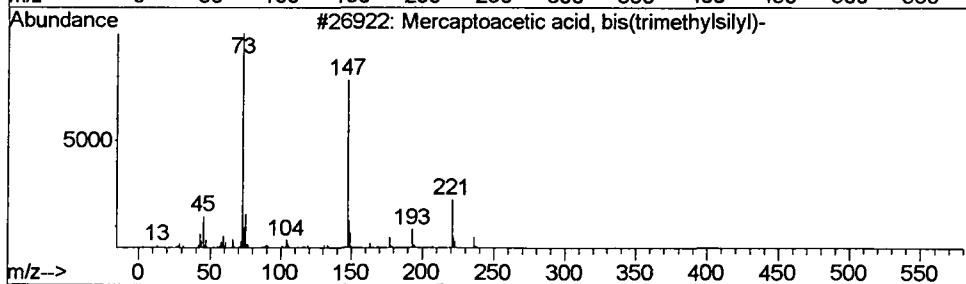
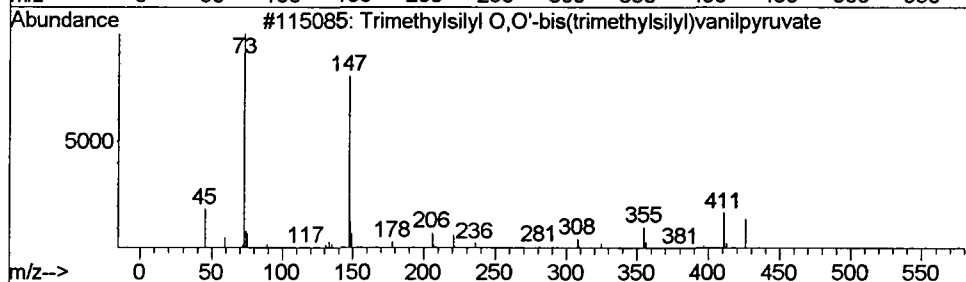
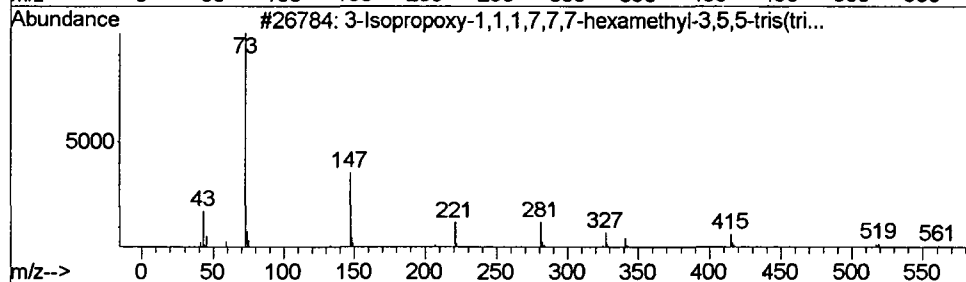
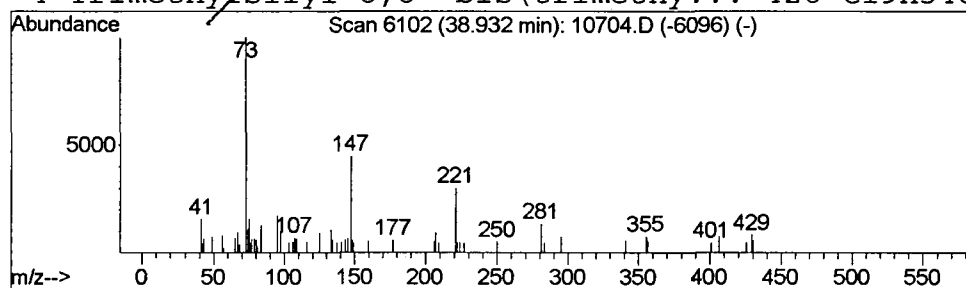
Vial: 15  
 Operator: Nefliu  
 Inst : Instrumen  
 Multiplr: 1.00

Quant Method : C:\MSDCHEM\1\METHODS\SCAN020507.M (RTE Integrator)  
 Title : Method 508 GC/MS Scan  
 Library : C:\DATABASE\nist98.1

\*\*\*\*\*  
 Peak Number 11 3-Isopropoxy-1,1,1,7,7,7-he... Concentration Rank 7

| R.T.  | EstConc   | Area  | Relative to ISTD | R.T.  |
|-------|-----------|-------|------------------|-------|
| 38.93 | 0.39 ug/l | 42673 | Chrysene-d12     | 39.94 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm     | CAS#        | Qual |
|------|----|---|-------------------------------------|-----|-------------|-------------|------|
| 1    |    |   | 3-Isopropoxy-1,1,1,7,7,7-hexamet... | 576 | C18H52O7Si7 | 071579-69-6 | 50   |
| 2    |    |   | Trimethylsilyl O,O'-bis(trimethy... | 426 | C19H34O5Si3 | 068595-71-1 | 43   |
| 3    |    |   | Mercaptoacetic acid, bis(trimeth... | 236 | C8H20O2SSi2 | 006398-62-5 | 40   |
| 4    |    |   | Trimethylsilyl O,O'-bis(trimethy... | 426 | C19H34O5Si3 | 068595-71-1 | 38   |



Operator ID: Nefliu Date Acquired: 14 May 2002 2:23 am  
 Data File: C:\MSDCHEM\1\DATA\020513\10704.D  
 Name: sample  
 Misc:  
 Method: C:\MSDCHEM\1\METHODS\SCAN020507.M (RTE Integrator)  
 Title: Method 508 GC/MS Scan  
 Library Searched: C:\DATABASE\nist98.1

| TIC Top Hit name                | RT    | EstConc | Units | Response | --Internal Standard-- |       |         |      |
|---------------------------------|-------|---------|-------|----------|-----------------------|-------|---------|------|
|                                 |       |         |       |          | #                     | RT    | Resp    | Conc |
| <del>Acetonitrile, dic...</del> | 3.66  | 0.6     | ug/l  | 54737    | 1                     | 11.38 | 3888190 | 40.0 |
| <del>2-Pentanone, 4-hy...</del> | 6.16  | 14.1    | ug/l  | 1371280  | 1                     | 11.38 | 3888190 | 40.0 |
| <del>Heptasiloxane, 1,...</del> | 28.36 | 0.4     | ug/l  | 54873    | 4                     | 31.72 | 5356370 | 40.0 |
| <del>Anthracene, 2-met...</del> | 31.09 | 0.3     | ug/l  | 46709    | 4                     | 31.72 | 5356370 | 40.0 |
| <del>Anthracene-d10-</del>      | 31.96 | 0.4     | ug/l  | 48682    | 4                     | 31.72 | 5356370 | 40.0 |
| <del>1,1,1,5,7,7,7-Hep...</del> | 32.20 | 0.4     | ug/l  | 53976    | 4                     | 31.72 | 5356370 | 40.0 |
| <del>Octasiloxane, 1,1...</del> | 34.55 | 0.3     | ug/l  | 40107    | 4                     | 31.72 | 5356370 | 40.0 |
| <del>3-Isopropoxy-1,1,...</del> | 36.28 | 0.4     | ug/l  | 43002    | 5                     | 39.94 | 4355360 | 40.0 |
| <del>t-Butyldimethyl(2...</del> | 37.70 | 0.4     | ug/l  | 42649    | 5                     | 39.94 | 4355360 | 40.0 |
| <del>Phosphonic acid, ...</del> | 38.54 | 5.9     | ug/l  | 647211   | 5                     | 39.94 | 4355360 | 40.0 |
| <del>3-Isopropoxy-1,1,...</del> | 38.93 | 0.4     | ug/l  | 42673    | 5                     | 39.94 | 4355360 | 40.0 |