

User: Vinci, David L  
 Westchester Dept. of Labs & Research  
 Date: 04/09/2002 Time: 09:46:38

Sample: AE07628  
 Analysis: \$GCMS GCMS Scan For Unknowns  
 Units: ug  
 Started: 4/9/02 09:46 Ended: 4/9/02 09:46 Analyst: DLV  
 Page: 1

|   | Component Name          | Viol | Result | Qualifier | MDL |
|---|-------------------------|------|--------|-----------|-----|
| 1 | Other unknown compounds |      | .      |           |     |
| 2 | Surr_2,4-DDD            |      | 89     |           | 5.0 |

**Comments for Sample AE07628 - Analysis \$GCMS**

Westchester Dept. of Labs & Research

User: Vinci, David L

Date: 04-09-2002 Time: 09:46:58

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

## Quantitation Report (QT Reviewed)

Data File : C:\MSDCHEM\1\DATA\020403\07628.D  
 Acq On : 4 Apr 2002 1:43 am  
 Sample : sample  
 Misc :  
 MS Integration Params: EVENTS.E  
 Quant Time: Apr 04 08:15:16 2002

Vial: 19  
 Operator: Nefliu  
 Inst : Instrumen  
 Multiplr: 1.00

Quant Results File: SCAN020403.RE

Quant Method : C:\MSDCHEM\1\METHODS\SCAN020403.M (Chemstation Integrator)  
 Title : Method 508 GC/MS Scan  
 Last Update : Thu Apr 04 08:10:12 2002  
 Response via : Initial Calibration  
 DataAcq Meth : 508SCAN1

| Internal Standards        | R.T.  | QIon | Response | Conc  | Units | Dev (Min) |
|---------------------------|-------|------|----------|-------|-------|-----------|
| 1) 1,4-Dichlorobenzene-d4 | 10.66 | 152  | 2543269  | 20.00 | ug/l  | 0.00      |
| 10) Naphthalene-d8        | 15.77 | 136  | 9276148  | 20.00 | ug/l  | 0.00      |
| 17) Acenaphthene-d10      | 23.66 | 164  | 5338604  | 20.00 | ug/l  | -0.01     |
| 31) Phenanthrene-d10      | 30.89 | 188  | 8695779  | 20.00 | ug/l  | 0.00      |
| 39) Chrysene-d12          | 39.09 | 240  | 6700530  | 20.00 | ug/l  | -0.02     |
| 45) Perylene-d12          | 42.30 | 264  | 3456595  | 20.00 | ug/l  | -0.01     |

## System Monitoring Compounds

|                  |       |     |          |      |        |      |
|------------------|-------|-----|----------|------|--------|------|
| 28) TCMX         | 26.72 | 207 | 186265   | 3.57 | ug/l   | 0.00 |
| Spiked Amount    |       |     | Recovery | =    | 71.40% |      |
| Target Compounds |       |     |          |      | 89     |      |
|                  |       |     |          |      | Qvalue |      |

-----  
 (#) = qualifier out of range (m) = manual integration (+) = signals summed  
 07628.D SCAN020403.M Thu Apr 04 10:36:13 2002 Page 1

## Quantitation Report (QT Reviewed)

Data File : C:\MSDCHEM\1\DATA\020403\07628.D

Vial: 19

Acq On : 4 Apr 2002 1:43 am

Operator: Nefliu

Sample : sample

Inst : Instrumen

Misc :

Multiplr: 1.00

MS Integration Params: EVENTS.E

Quant Time: Apr 4 10:36 2002

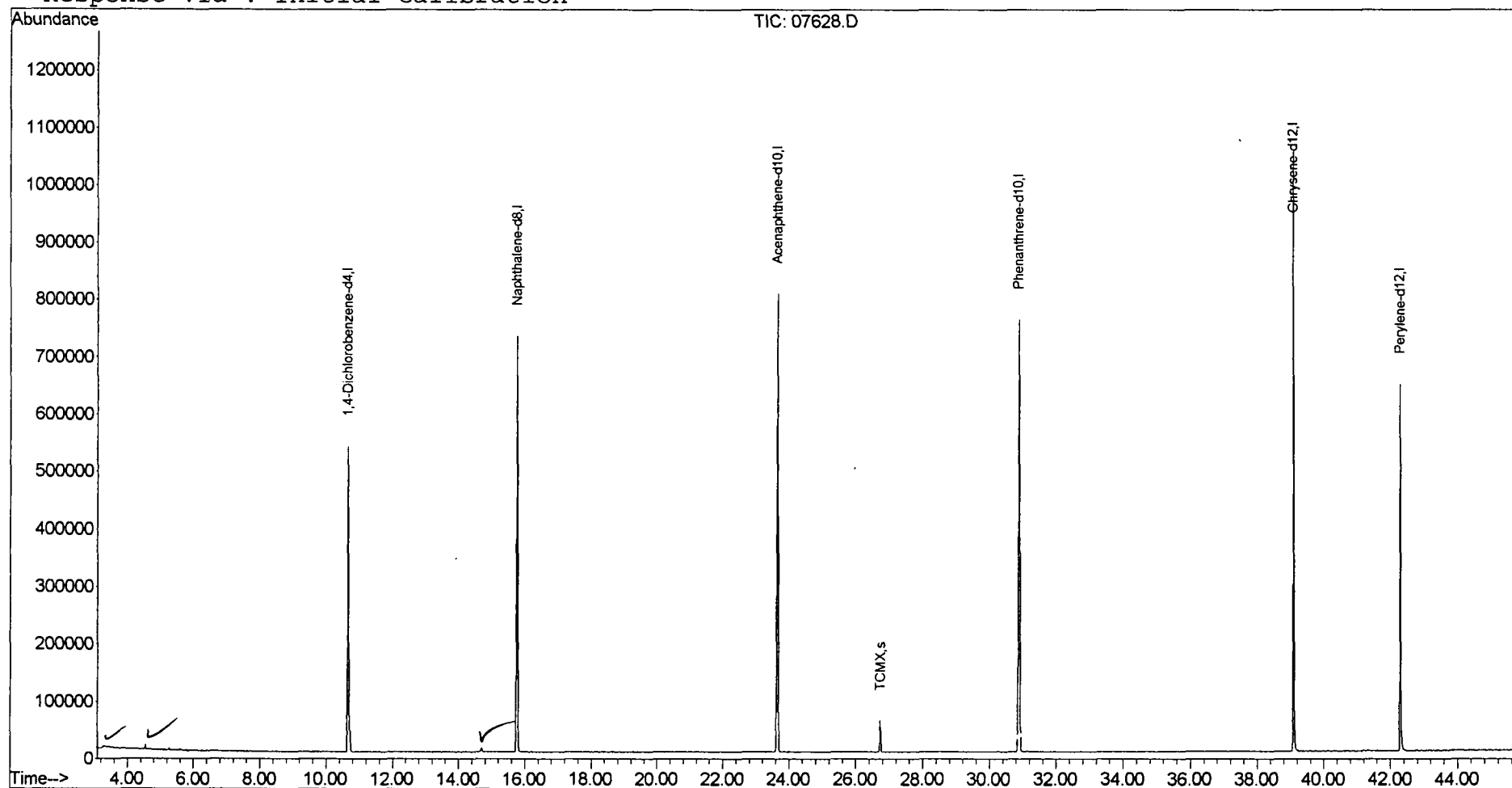
Quant Results File: SCAN020403.RES

Method : C:\MSDCHEM\1\METHODS\SCAN020403.M (Chemstation Integrator)

Title : Method 508 GC/MS Scan

Last Update : Thu Apr 04 08:40:00 2002

Response via : Initial Calibration



## LSC Area Percent Report

Data File : C:\MSDCHEM\1\DATA\020403\07628.D  
Acq On : 4 Apr 2002 1:43 am  
Sample : sample  
Misc :  
MS Integration Params: LSCINT.e

Vial: 19  
Operator: Nefliu  
Inst : Instrumen  
Multiplr: 1.00

Method : C:\MSDCHEM\1\METHODS\SCAN020403.M (Chemstation Integrator)  
Title : Method 508 GC/MS Scan

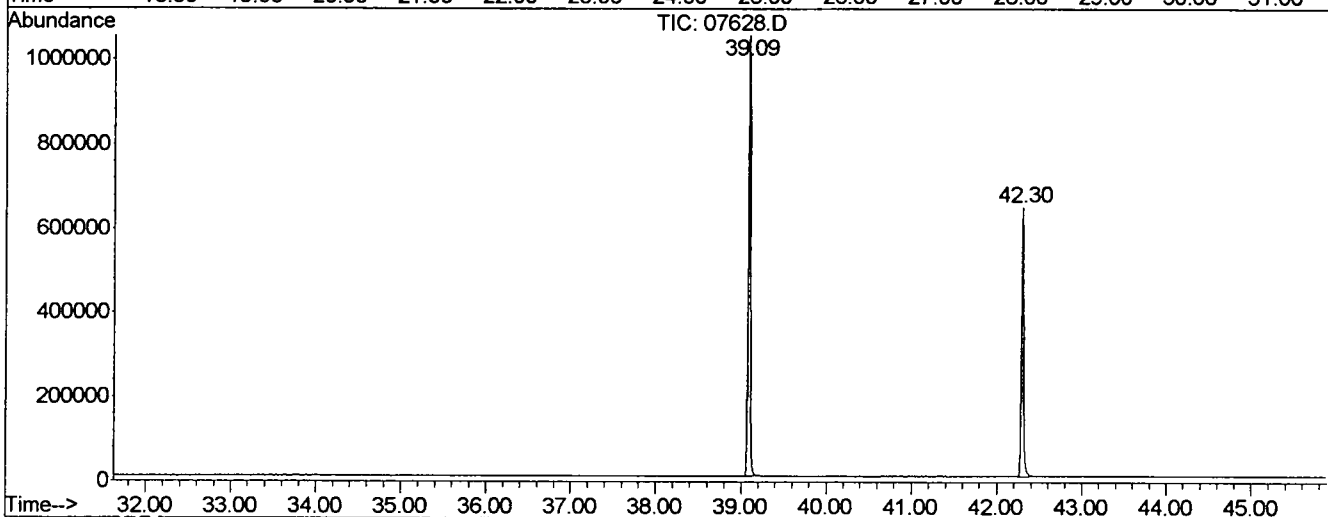
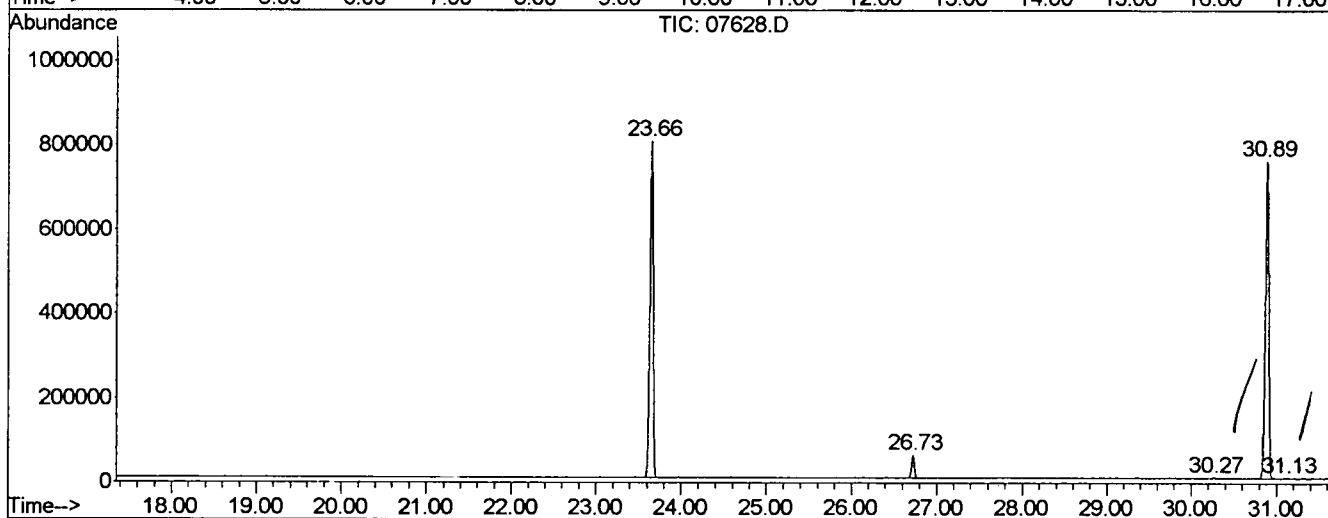
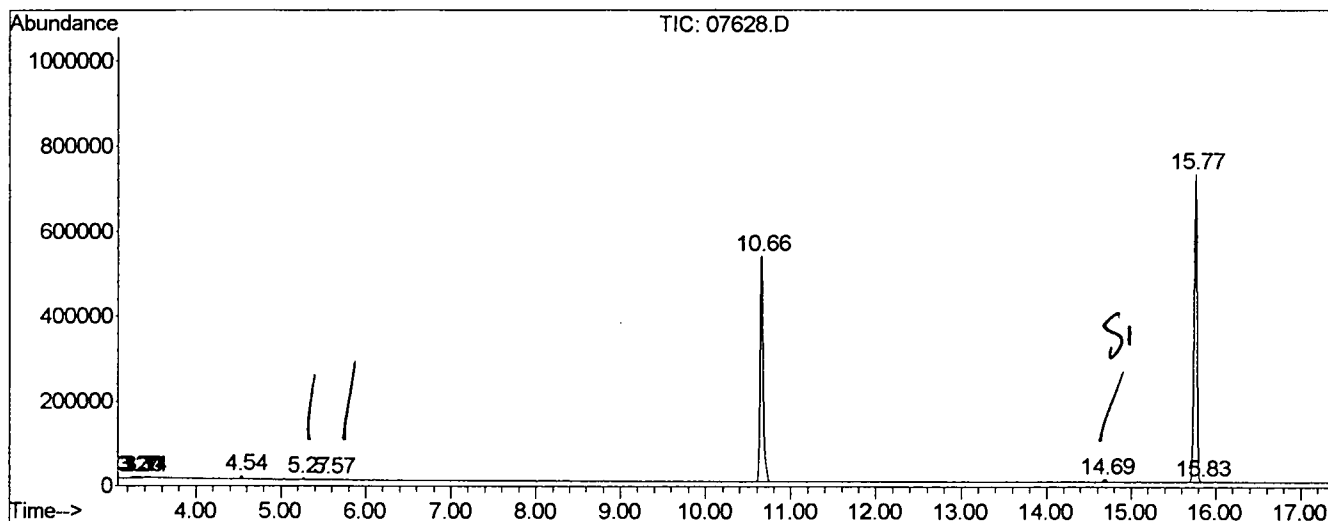
Signal : TIC

| 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         | 3.267       | 21            | 32          | 34           | PV 3     | 3681           | 92108         | 0.43%           | 0.088%        |
| 2         | 3.297       | 34            | 37          | 39           | VV 2     | 4028           | 63174         | 0.30%           | 0.061%        |
| 3         | 3.314       | 39            | 40          | 43           | VV 2     | 3988           | 63538         | 0.30%           | 0.061%        |
| 4         | 3.344       | 43            | 45          | 51           | VV 3     | 3765           | 90339         | 0.42%           | 0.087%        |
| 5         | 4.536       | 242           | 248         | 256          | VV 2     | 8351           | 138260        | 0.65%           | 0.133%        |
| 6         | 5.271       | 362           | 373         | 377          | VV 4     | 4402           | 65693         | 0.31%           | 0.063%        |
| 7         | 5.570       | 421           | 424         | 428          | PV 2     | 2017           | 27062         | 0.13%           | 0.026%        |
| 8         | 10.659      | 1279          | 1290        | 1309         | PV       | 525040         | 13205433      | 62.11%          | 12.665%       |
| 9         | 14.689      | 1967          | 1976        | 1987         | VV 3     | 6764           | 180235        | 0.85%           | 0.173%        |
| 10        | 15.764      | 2146          | 2159        | 2169         | VV       | 715641         | 17441621      | 82.04%          | 16.728%       |
| 11        | 15.829      | 2169          | 2170        | 2178         | VV 2     | 2178           | 40555         | 0.19%           | 0.039%        |
| 12        | 23.655      | 3477          | 3502        | 3517         | PV       | 782993         | 20593223      | 96.86%          | 19.751%       |
| 13        | 26.722      | 4016          | 4024        | 4033         | VV 2     | 52805          | 1171907       | 5.51%           | 1.124%        |
| 14        | 30.271      | 4621          | 4628        | 4638         | VV 2     | 2378           | 62833         | 0.30%           | 0.060%        |
| 15        | 30.894      | 4713          | 4734        | 4751         | PV 2     | 750877         | 21260418      | 100.00%         | 20.391%       |
| 16        | 31.129      | 4765          | 4774        | 4780         | PV 2     | 1887           | 57212         | 0.27%           | 0.055%        |
| 17        | 39.091      | 6120          | 6129        | 6147         | PV 2     | 1025466        | 18479666      | 86.92%          | 17.724%       |
| 18        | 42.299      | 6665          | 6675        | 6704         | PV       | 636156         | 11231020      | 52.83%          | 10.772%       |

Sum of corrected areas: 104264299

# LSC Report - Integrated Chromatogram

File : C:\MSDCHEM\1\DATA\020403\07628.D  
 Operator : Nefliu  
 Acquired : 4 Apr 2002 1:43 am using AcqMethod 508SCAN1  
 Instrument : Instrumen  
 Sample Name: sample  
 Misc Info :  
 Vial Number: 19  
 Quant File :SCAN020403.RES (Chemstation Integrator)



## Library Search Compound Report

Data File : C:\MSDCHEM\1\DATA\020403\07628.D  
Acq On : 4 Apr 2002 1:43 am  
Sample : sample  
Misc :  
MS Integration Params: LSCINT.e

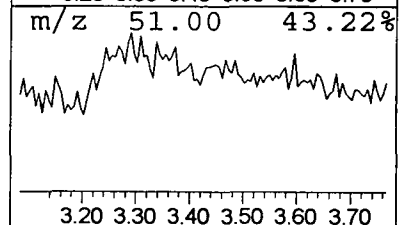
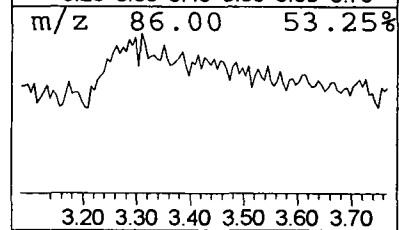
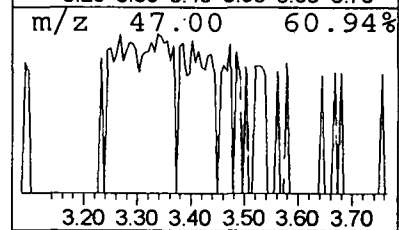
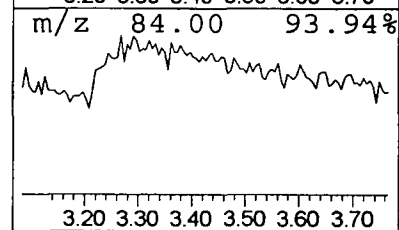
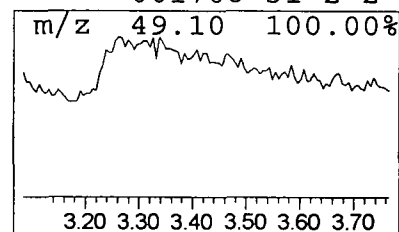
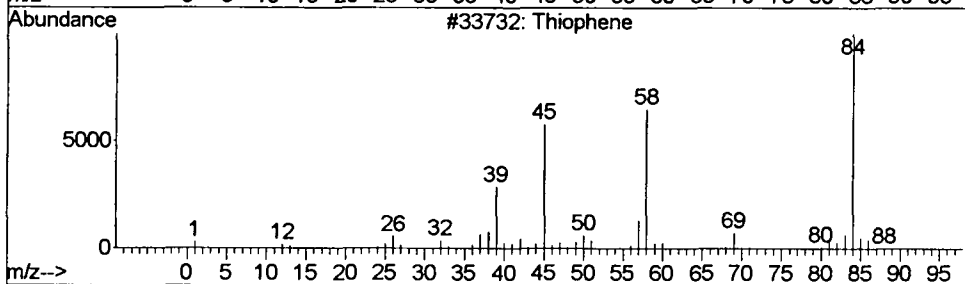
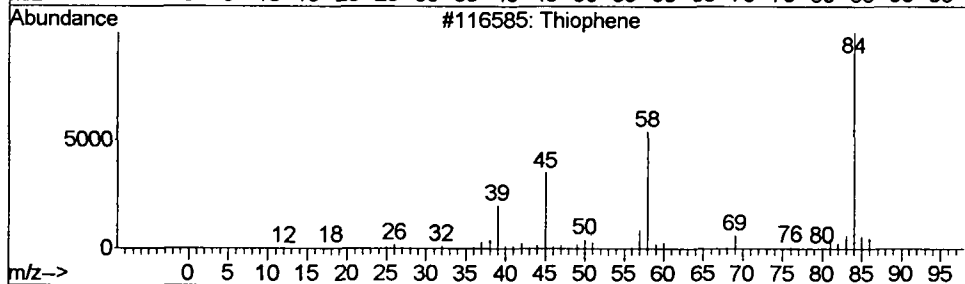
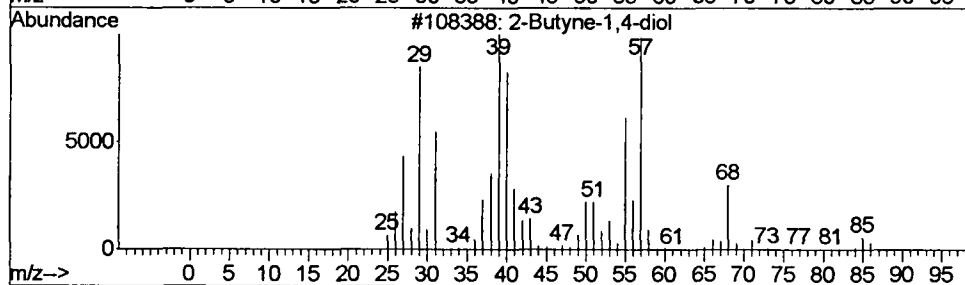
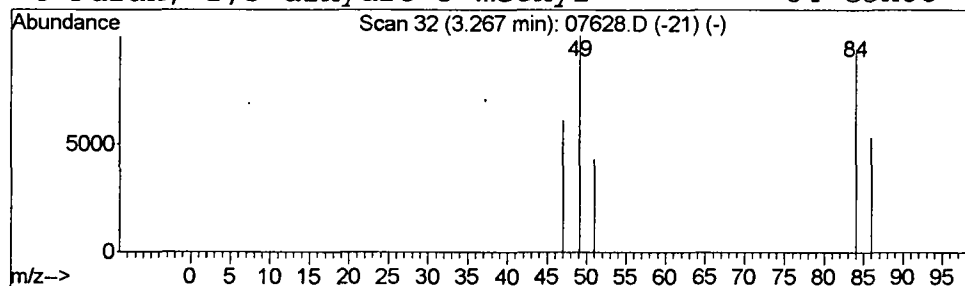
Vial: 19  
Operator: Nefliu  
Inst : Instrumen  
Multiplr: 1.00

Quant Method : C:\MSDCHEM\1\METHODS\SCAN020403.M (Chemstation Integrator)  
Title : Method 508 GC/MS Scan  
Library : C:\DATABASE\nist98.1

\*\*\*\*\*  
Peak Number 1 2-Butyne-1,4-diol Concentration Rank 3

| R.T. | EstConc   | Area  | Relative to ISTD       | R.T.  |
|------|-----------|-------|------------------------|-------|
| 3.27 | 0.14 ug/l | 92108 | 1,4-Dichlorobenzene-d4 | 10.66 |

| Hit# | of | 5 | Tentative ID                 | MW | MolForm | CAS#        | Qual |
|------|----|---|------------------------------|----|---------|-------------|------|
| 1    |    |   | 2-Butyne-1,4-diol            | 86 | C4H6O2  | 000110-65-6 | 3    |
| 2    |    |   | Thiophene                    | 84 | C4H4S   | 000110-02-1 | 2    |
| 3    |    |   | Thiophene                    | 84 | C4H4S   | 000110-02-1 | 2    |
| 4    |    |   | Furan, 2,5-dihydro-3-methyl- | 84 | C5H8O   | 001708-31-2 | 2    |



# Library Search Compound Report

Data File : C:\MSDCHEM\1\DATA\020403\07628.D  
 Acq On : 4 Apr 2002 1:43 am  
 Sample : sample  
 Misc :  
 MS Integration Params: LSCINT.e

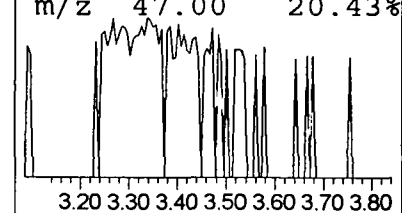
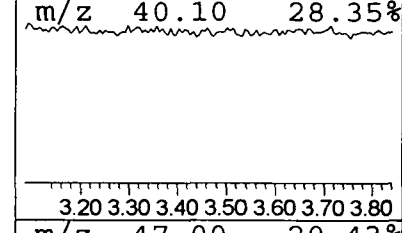
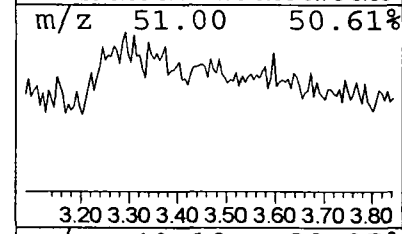
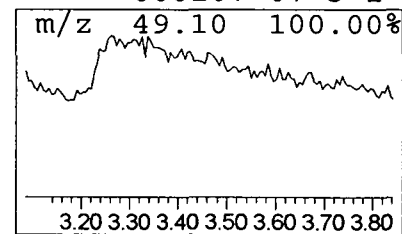
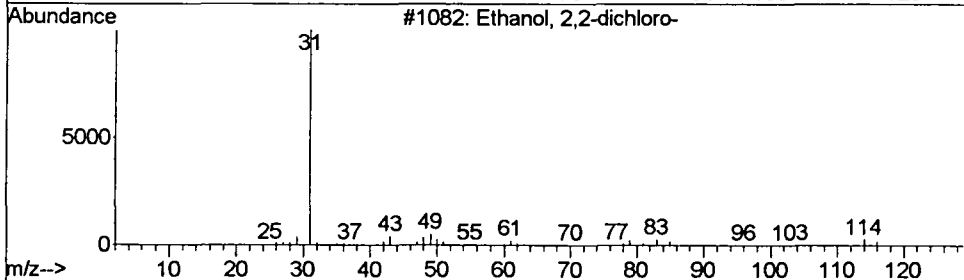
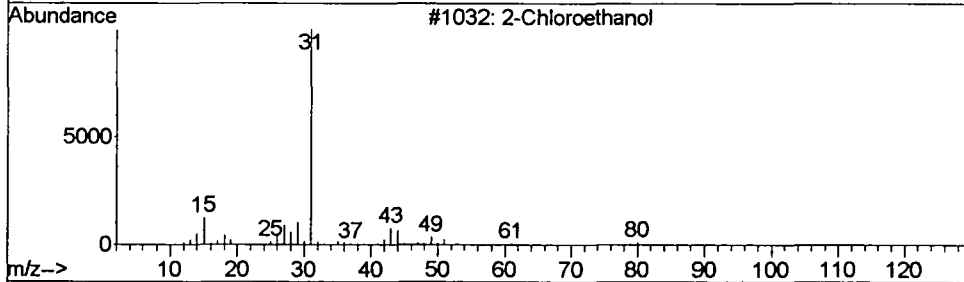
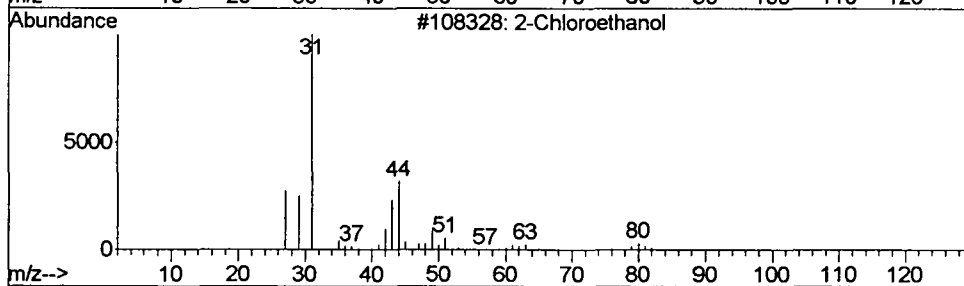
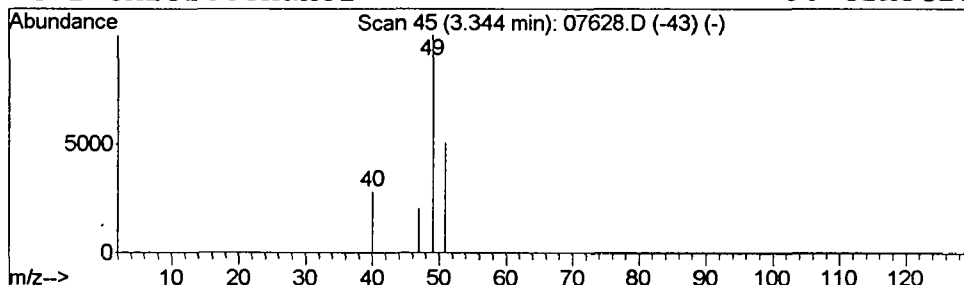
Vial: 19  
 Operator: Nefliu  
 Inst : Instrumen  
 Multiplr: 1.00

Quant Method : C:\MSDCHEM\1\METHODS\SCAN020403.M (Chemstation Integrator)  
 Title : Method 508 GC/MS Scan  
 Library : C:\DATABASE\nist98.1

\*\*\*\*\*  
 Peak Number 2 2-Chloroethanol Concentration Rank 4

| R.T. | EstConc   | Area  | Relative to ISTD       | R.T.  |
|------|-----------|-------|------------------------|-------|
| 3.34 | 0.14 ug/l | 90339 | 1,4-Dichlorobenzene-d4 | 10.66 |

| Hit# | of | 5 | Tentative ID           | MW  | MolForm  | CAS#        | Qual |
|------|----|---|------------------------|-----|----------|-------------|------|
| 1    |    |   | 2-Chloroethanol        | 80  | C2H5ClO  | 000107-07-3 | 2    |
| 2    |    |   | 2-Chloroethanol        | 80  | C2H5ClO  | 000107-07-3 | 2    |
| 3    |    |   | Ethanol, 2,2-dichloro- | 114 | C2H4Cl2O | 000598-38-9 | 1    |
| 4    |    |   | 2-Chloroethanol        | 80  | C2H5ClO  | 000107-07-3 | 1    |



# Library Search Compound Report

Data File : C:\MSDCHEM\1\DATA\020403\07628.D  
 Acq On : 4 Apr 2002 1:43 am  
 Sample : sample  
 Misc :  
 MS Integration Params: LSCINT.e

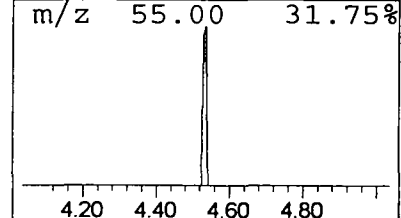
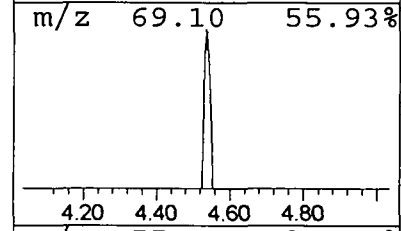
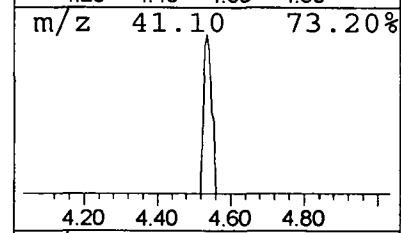
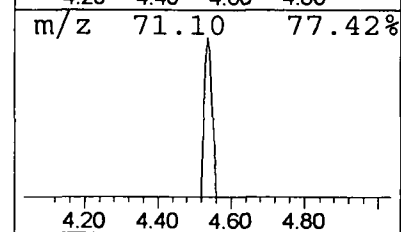
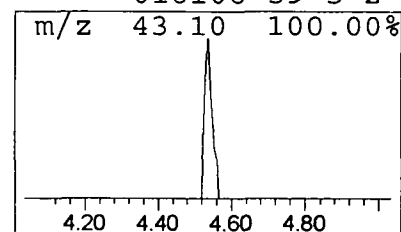
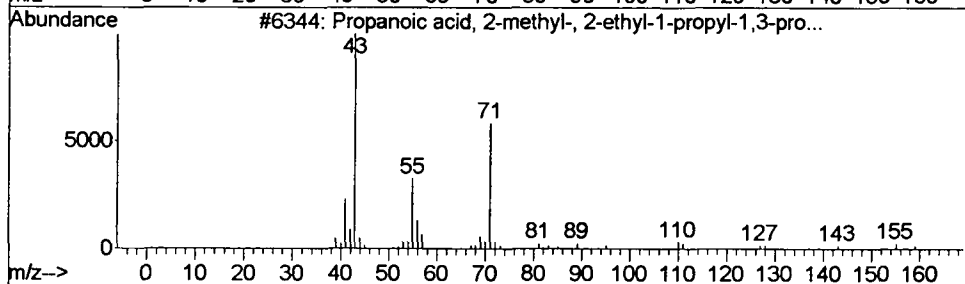
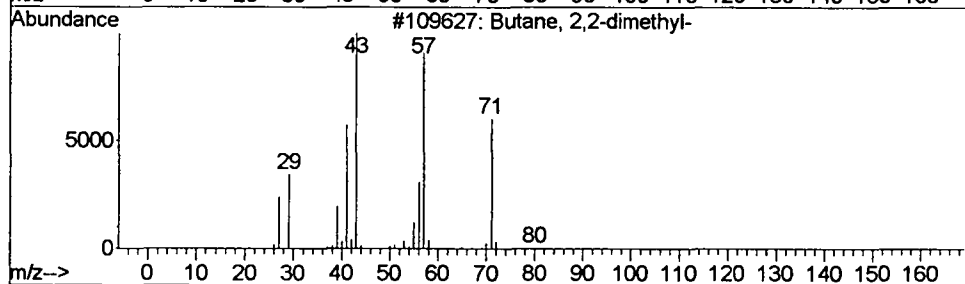
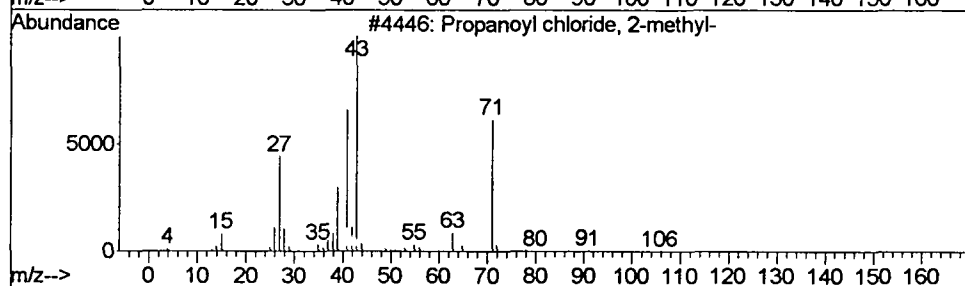
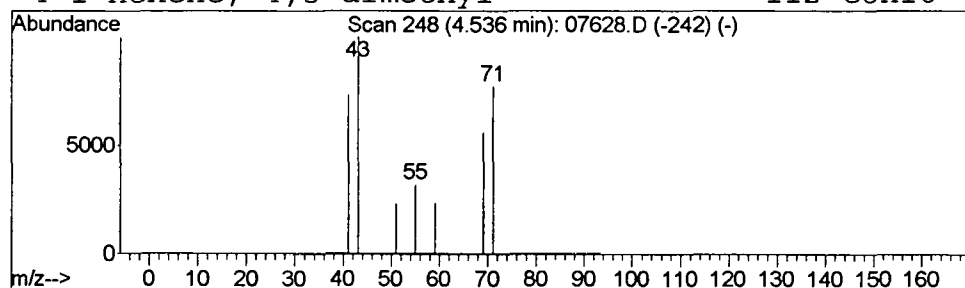
Vial: 19  
 Operator: Nefliu  
 Inst : Instrumen  
 Multiplr: 1.00

Quant Method : C:\MSDCHEM\1\METHODS\SCAN020403.M (Chemstation Integrator)  
 Title : Method 508 GC/MS Scan  
 Library : C:\DATABASE\nist98.1

\*\*\*\*\*  
 Peak Number 3 Propanoyl chloride, 2-methyl- Concentration Rank 1

| R.T. | EstConc   | Area   | Relative to ISTD       | R.T.  |
|------|-----------|--------|------------------------|-------|
| 4.54 | 0.21 ug/l | 138260 | 1,4-Dichlorobenzene-d4 | 10.66 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm  | CAS#        | Qual |
|------|----|---|-------------------------------------|-----|----------|-------------|------|
| 1    |    |   | Propanoyl chloride, 2-methyl-       | 106 | C4H7ClO  | 000079-30-1 | 2    |
| 2    |    |   | Butane, 2,2-dimethyl-               | 86  | C6H14    | 000075-83-2 | 2    |
| 3    |    |   | Propanoic acid, 2-methyl-, 2-eth... | 286 | C16H30O4 | 074367-30-9 | 2    |
| 4    |    |   | 1-Hexene, 4,5-dimethyl-             | 112 | C8H16    | 016106-59-5 | 2    |



## Library Search Compound Report

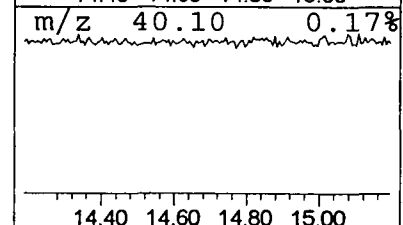
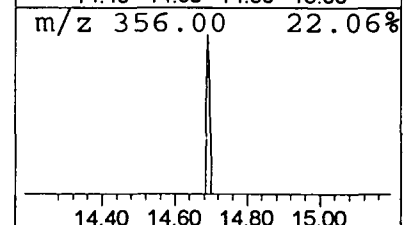
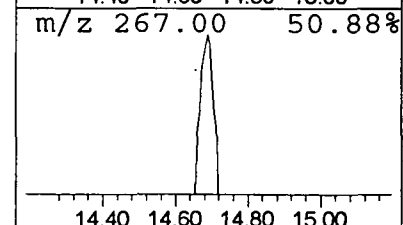
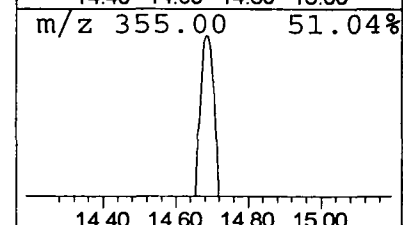
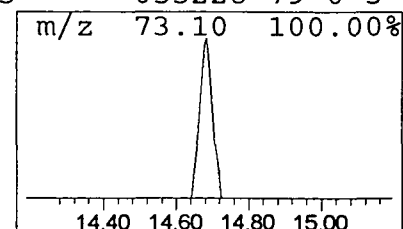
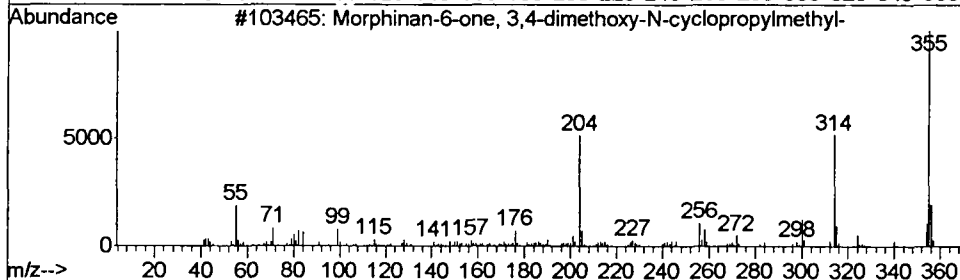
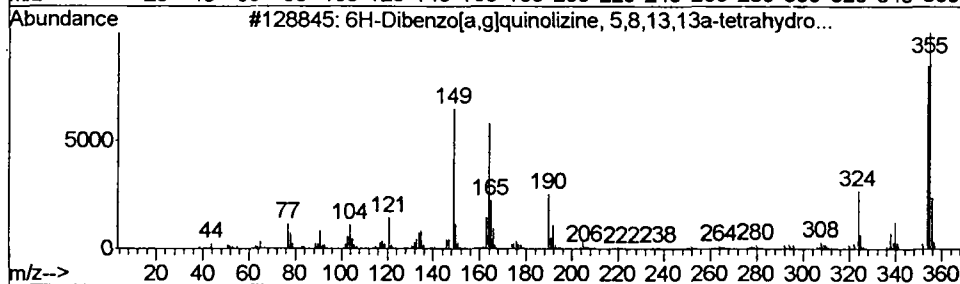
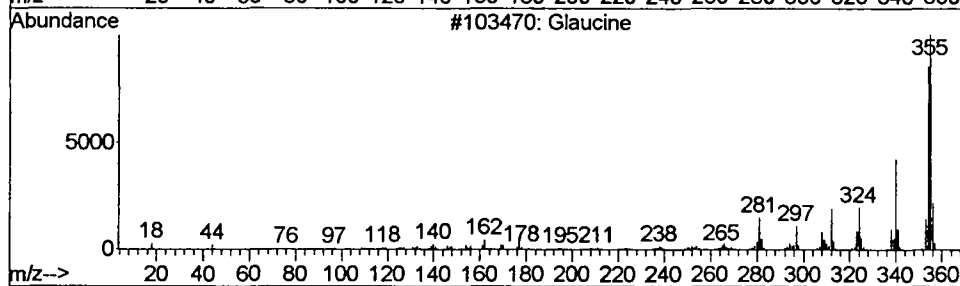
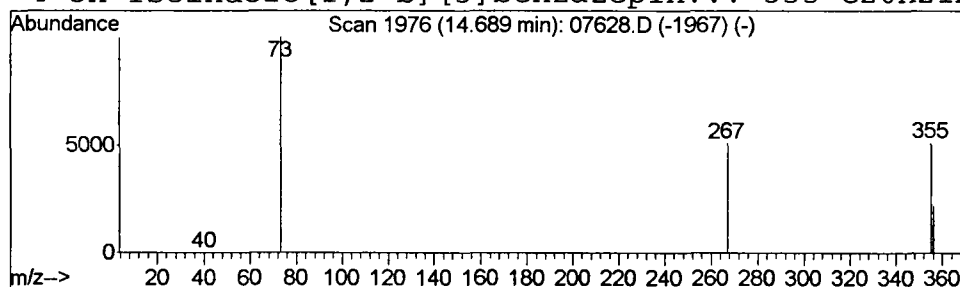
Data File : C:\MSDCHEM\1\DATA\020403\07628.D  
Acq On : 4 Apr 2002 1:43 am  
Sample : sample  
Misc :  
MS Integration Params: LSCINT.e

Vial: 19  
Operator: Nefliu  
Inst : Instrumen  
Multiplr: 1.00

Quant Method : C:\MSDCHEM\1\METHODS\SCAN020403.M (Chemstation Integrator)  
Title : Method 508 GC/MS Scan  
Library : C:\DATABASE\nist98.1

\*\*\*\*\*  
Peak Number 4 Glaucine Concentration Rank 2

| R.T.    | EstConc                             | Area         | Relative to ISTD | R.T.      |              |      |
|---------|-------------------------------------|--------------|------------------|-----------|--------------|------|
| 14.69   | 0.21 ug/l                           | 180235       | Naphthalene-d8   | 15.77     |              |      |
| Hit# of | 5                                   | Tentative ID | MW               | MolForm   | CAS#         | Qual |
| 1       | Glaucine                            |              | 355              | C21H25NO4 | 000475-81-0  | 3    |
| 2       | 6H-Dibenzo[a,g]quinolizine, 5,8,... |              | 355              | C21H25NO4 | 002934-97-6  | 3    |
| 3       | Morphinan-6-one, 3,4-dimethoxy-N... |              | 355              | C22H29NO3 | 1000126-19-3 | 3    |
| 4       | 5H-Isoindolo[1,2-b][3]benzazepin... |              | 355              | C20H21NO5 | 055228-79-0  | 3    |



## Tentatively Identified Compound (LSC) summary

Operator ID: Nefliu Date Acquired: 4 Apr 2002 1:43 am  
Data File: C:\MSDCHEM\1\DATA\020403\07628.D  
Name: sample  
Misc:  
Method: C:\MSDCHEM\1\METHODS\SCAN020403.M (Chemstation Integrator)  
Title: Method 508 GC/MS Scan  
Library Searched: C:\DATABASE\nist98.l

| TIC Top Hit name     | RT    | EstConc | Units | Response | # | RT    | Resp     | Conc |
|----------------------|-------|---------|-------|----------|---|-------|----------|------|
| 2-Butyne-1,4-diol    | 3.27  | 0.1     | ug/l  | 92108    | 1 | 10.66 | 13205400 | 20.0 |
| 2-Chloroethanol      | 3.34  | 0.1     | ug/l  | 90339    | 1 | 10.66 | 13205400 | 20.0 |
| Propanoyl chlorid... | 4.54  | 0.2     | ug/l  | 138260   | 1 | 10.66 | 13205400 | 20.0 |
| Glaucine             | 14.69 | 0.2     | ug/l  | 180235   | 2 | 15.77 | 17441600 | 20.0 |