

## Quantitation Report

(QT Reviewed)

Data File : C:\MSDCHEM\1\DATA\022702\02577.D

Acq On : 28 Feb 2002 1:29 am

Sample : GC SCAN 2/4 B

Misc :

MS Integration Params: BENZO.P

Quant Time: Mar 01 09:03:20 2002

Vial: 20

Operator: McLean

Inst : Instrumen

Multiplr: 1.00

Quant Results File: 5080202.RES

Quant Method : C:\MSDCHEM\1\5973N\5080202.M (RTE Integrator)

Title : Quantitation For Method 508gcscan

Last Update : Mon Feb 25 19:17:14 2002

Response via : Initial Calibration

DataAcq Meth : 508SCAN

| Internal Standards        | R.T.  | QIon | Response | Conc  | Units | Dev(Min) |
|---------------------------|-------|------|----------|-------|-------|----------|
| 1) 1,4-Dichlorobenzene-d4 | 9.58  | 150  | 2315695  | 20.00 | ug/L  | 0.00     |
| 10) Naphthalene-d8        | 13.06 | 136  | 6024409  | 20.00 | ug/L  | 0.00     |
| 17) Acenaphthene-d10      | 18.18 | 164  | 3335305  | 20.00 | ug/L  | 0.00     |
| 29) Phenanthrene-d10      | 22.54 | 188  | 5191389  | 20.00 | ug/L  | 0.00     |
| 38) Chrysene-d12          | 30.35 | 240  | 4865611  | 20.00 | ug/L  | 0.00     |
| 44) Perylene-d12          | 34.23 | 264  | 3482972  | 20.00 | ug/L  | 0.00     |

## System Monitoring Compounds

|                  |       |     |          |      |        |      |
|------------------|-------|-----|----------|------|--------|------|
| 37) 2,4-ddd surr | 27.46 | 235 | 219514   | 2.66 | ug/L   | 0.00 |
| Spiked Amount    | 5.000 |     | Recovery | =    | 53.20% |      |

## Target Compounds

|                                |       |     |       |      |      | Qvalue |
|--------------------------------|-------|-----|-------|------|------|--------|
| 2) N-nitroso-dimethyl amine    | 0.00  | 74  | 0     | N.D. | d    |        |
| 3) bis(2-chloroethyl) ether    | 0.00  | 93  | 0     | N.D. |      |        |
| 4) 1,3 - Dichlorobenze         | 0.00  | 146 | 0     | N.D. |      |        |
| 5) 1,4 - Dichlorobenzene       | 0.00  | 146 | 0     | N.D. |      |        |
| 6) 1,2 - Dichlorobenzene       | 0.00  | 146 | 0     | N.D. |      |        |
| 7) 2,2' - oxybis (1-Chloropr   | 0.00  | 45  | 0     | N.D. |      |        |
| 8) N-Nitroso-di-n-propylamine  | 0.00  | 70  | 0     | N.D. |      |        |
| 9) Hexachloroethane            | 0.00  | 117 | 0     | N.D. |      |        |
| 11) Nitrobenzene               | 0.00  | 77  | 0     | N.D. |      |        |
| 12) Isophorone                 | 0.00  | 82  | 0     | N.D. |      |        |
| 13) bis(2-Chloroethoxy) methan | 0.00  | 93  | 0     | N.D. |      |        |
| 14) 1,2,4-Trichlorobenzene     | 0.00  | 180 | 0     | N.D. |      |        |
| 15) Naphthalene                | 0.00  | 128 | 0     | N.D. |      |        |
| 16) Hexachlorobutadiene        | 0.00  | 225 | 0     | N.D. |      |        |
| 18) Hexachlorocyclopentadiene  | 0.00  | 237 | 0     | N.D. |      |        |
| 19) 2-chloronaphthalene        | 0.00  | 162 | 0     | N.D. |      |        |
| 20) Dimethylphthalate          | 0.00  | 163 | 0     | N.D. |      |        |
| 21) 2,6-Dinitrotoluene         | 0.00  | 165 | 0     | N.D. | d    |        |
| 22) Acenaphthylene             | 0.00  | 152 | 0     | N.D. |      |        |
| 23) Acenaphthene               | 0.00  | 153 | 0     | N.D. |      |        |
| 24) 2,4-Dinitrotoluene         | 0.00  | 165 | 0     | N.D. |      |        |
| 25) Diethylphthalate           | 0.00  | 149 | 0     | N.D. |      |        |
| 26) 4-Chlorophenyl-phenyl ethe | 0.00  | 204 | 0     | N.D. |      |        |
| 27) Fluorene                   | 0.00  | 166 | 0     | N.D. |      |        |
| 28) Azobenzen/ 1,2-Diphenylhyd | 0.00  | 77  | 0     | N.D. |      |        |
| 30) N-Nitrosodiphenylamine     | 0.00  | 169 | 0     | N.D. |      |        |
| 31) 4-Bromophenyl-phenyl ether | 0.00  | 248 | 0     | N.D. |      |        |
| 32) Hexachlorobenzene          | 0.00  | 284 | 0     | N.D. |      |        |
| 33) Phenanthrene               | 0.00  | 178 | 0     | N.D. |      |        |
| 34) Anthracene                 | 0.00  | 178 | 0     | N.D. |      |        |
| 35) Di-n-butylphthalate        | 24.67 | 149 | 8069  | 0.02 | ug/L | 79     |
| 36) Fluoranthene               | 0.00  | 202 | 0     | N.D. |      |        |
| 39) Pyrene                     | 0.00  | 202 | 0     | N.D. |      |        |
| 40) Butylbenzylphthalate       | 0.00  | 149 | 0     | N.D. |      |        |
| 41) Benzo (a) anthracene       | 0.00  | 228 | 0     | N.D. | d    |        |
| 42) Bis (2-ethylhexyl) phthala | 30.95 | 149 | 19463 | 0.09 | ug/L | 91     |
| 43) Chrysene                   | 30.35 | 228 | 13109 | 0.04 | ug/L | 64     |
| 45) Di-n-Octylphthalate        | 0.00  | 149 | 0     | N.D. |      |        |
| 46) Benzo (b) fluoranthene     | 0.00  | 252 | 0     | N.D. |      |        |

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

02577.D 5080202.M Fri Mar 01 09:03:50 2002

## Quantitation Report (QT Reviewed)

Data File : C:\MSDCHEM\1\DATA\022702\02577.D  
Acq On : 28 Feb 2002 1:29 am  
Sample : GC SCAN 2/4 B  
Misc :

Vial: 20  
Operator: McLean  
Inst : Instrumen  
Multiplr: 1.00

MS Integration Params: BENZO.P

Quant Time: Mar 01 09:03:20 2002

Quant Results File: 5080202.RES

Quant Method : C:\MSDCHEM\1\5973N\5080202.M (RTE Integrator)

Title : Quantitation For Method 508gcscan

Last Update : Mon Feb 25 19:17:14 2002

Response via : Initial Calibration

DataAcq Meth : 508SCAN

| Compound                     | R.T. | QIon | Response | Conc | Unit | Qvalue |
|------------------------------|------|------|----------|------|------|--------|
| 47) Benzo (k) fluoranthene   | 0.00 | 252  | 0        | N.D. |      |        |
| 48) Benzo (a) pyrene         | 0.00 | 252  | 0        | N.D. | d    |        |
| 49) Indeno (1,2,3-cd) pyrene | 0.00 | 276  | 0        | N.D. |      |        |
| 50) Dibenz (a,h) anthracene  | 0.00 | 278  | 0        | N.D. |      |        |
| 51) Benzo (g,h,i) perylene   | 0.00 | 276  | 0        | N.D. |      |        |

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

02577.D 5080202.M Fri Mar 01 09:03:50 2002

Page 2

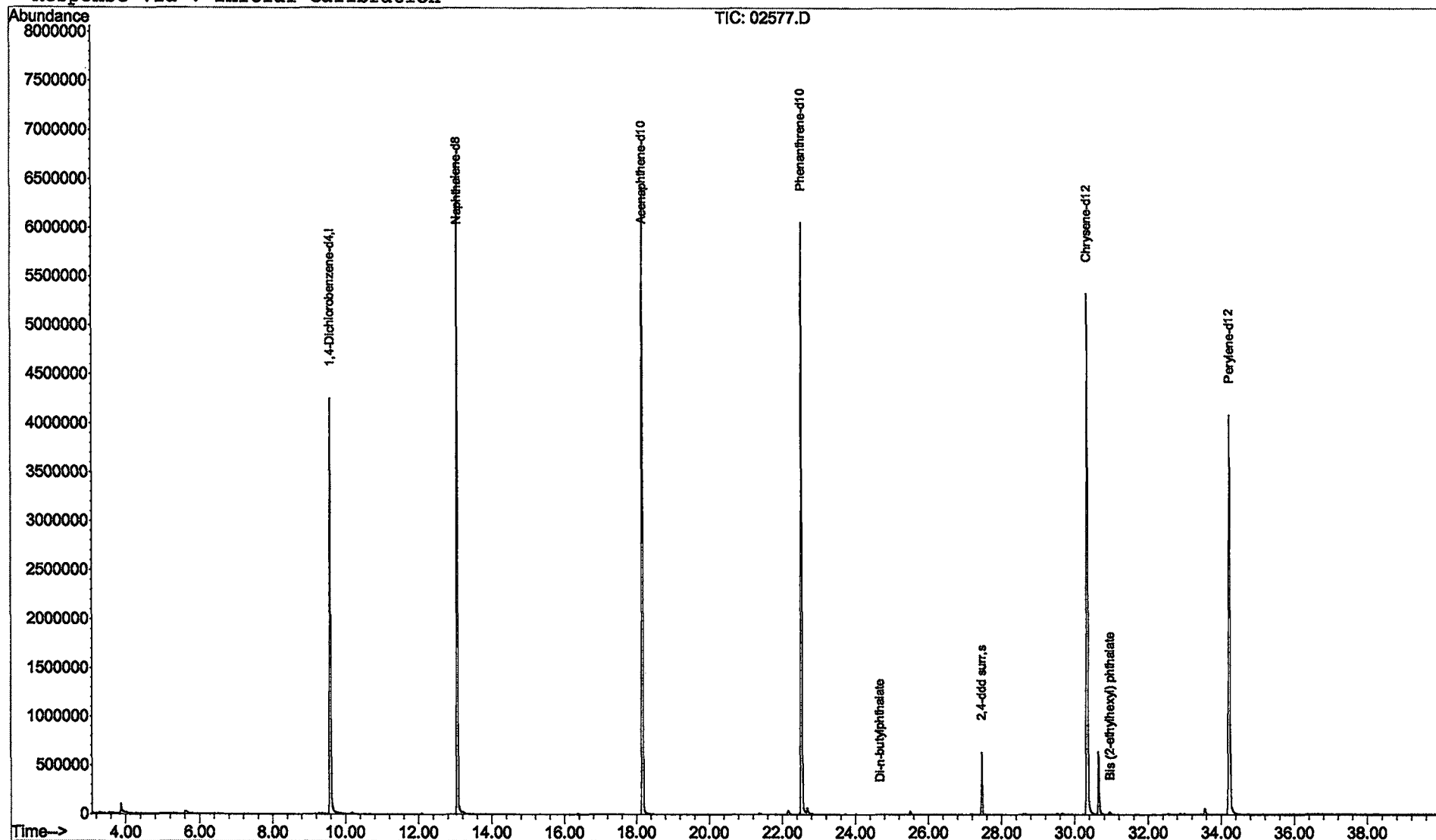
Quantitation Report (QI Reviewed)

Data File : C:\MSDCHEM\1\DATA\022702\02577.D  
 Acq On : 28 Feb 2002 1:29 am  
 Sample : GC SCAN 2/4 B  
 Misc :  
 MS Integration Params: BENZO.P  
 Quant Time: Mar 1 9:03 2002

Vial: 20  
 Operator: McLean  
 Inst : Instrumen  
 Multiplr: 1.00

Quant Results File: 5080202.RES

Method : C:\MSDCHEM\1\5973N\5080202.M (RTE Integrator)  
 Title : Quantitation For Method 508gcscan  
 Last Update : Mon Feb 25 19:17:14 2002  
 Response via : Initial Calibration



## LSC Area Percent Report

Data File : C:\MSDCHEM\1\DATA\022702\02577.D  
Acq On : 28 Feb 2002 1:29 am  
Sample : GC SCAN 2/4 B  
Misc :  
MS Integration Params: LSCINT.P

Vial: 20  
Operator: McLean  
Inst : Instrumen  
Multiplr: 1.00

Method : C:\MSDCHEM\1\5973N\5080202.M (RTE Integrator)  
Title : Quantitation For Method 508gcscan  
Smoothing : OFF Filtering: 5  
Sampling : 1 Min Area: 1 % of largest Peak  
Start Thrs: 0.2 Max Peaks: 100  
Stop Thrs : 0 Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >  
Peak separation: 5

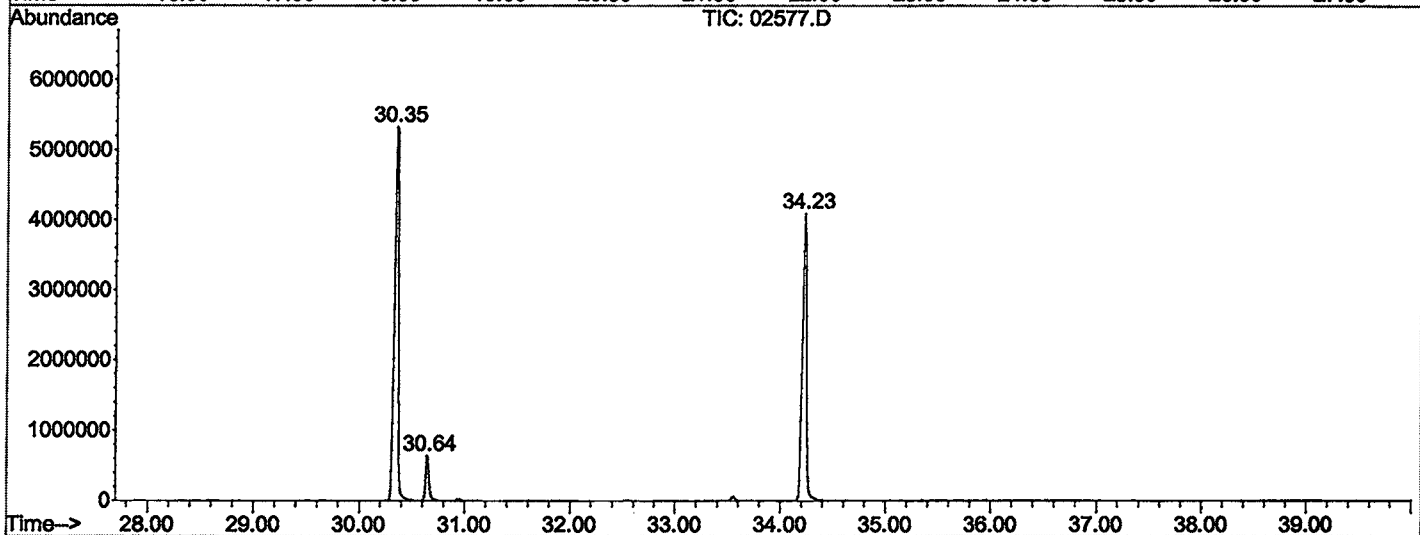
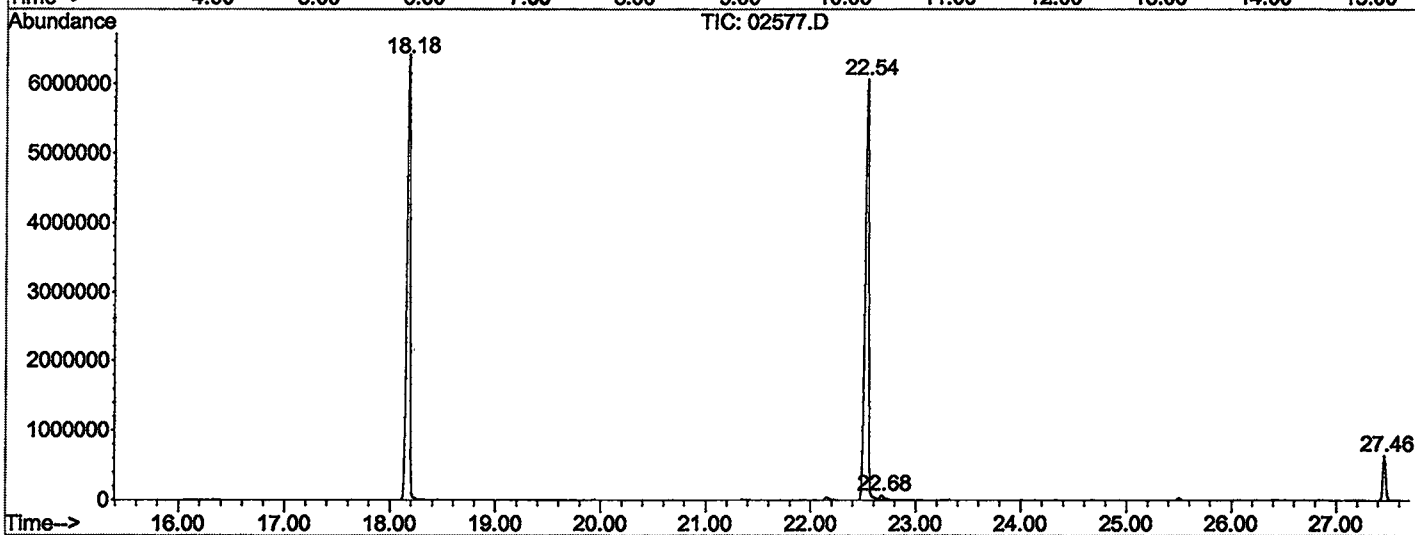
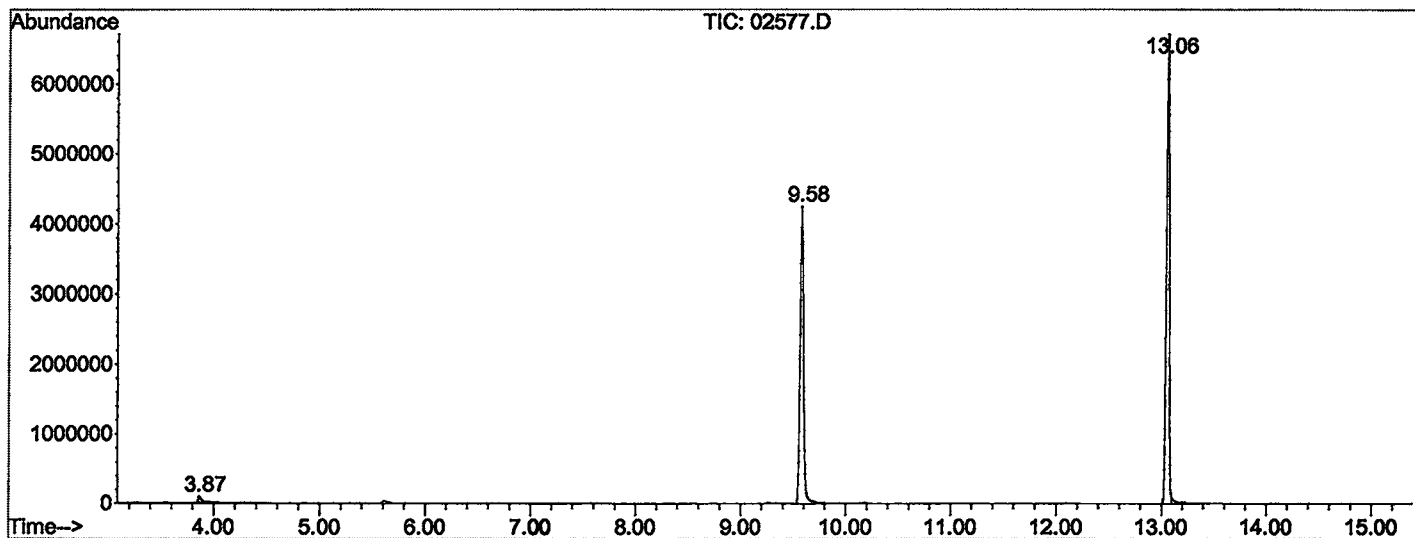
Signal : TIC

| 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         | 3.865       | 84            | 87          | 94           | rBV      | 92116          | 202285        | 1.41%           | 0.254%        |
| 2         | 9.579       | 711           | 717         | 736          | rBV      | 4253121        | 9729346       | 68.00%          | 12.238%       |
| 3         | 13.062      | 1094          | 1101        | 1105         | rBV      | 6713902        | 12914816      | 90.27%          | 16.244%       |
| 4         | 18.178      | 1657          | 1665        | 1679         | rBV      | 6424267        | 13912919      | 97.24%          | 17.500%       |
| 5         | 22.541      | 2138          | 2146        | 2158         | rBV2     | 6060995        | 14307515      | 100.00%         | 17.996%       |
| 6         | 22.677      | 2158          | 2161        | 2169         | rVB2     | 59445          | 148442        | 1.04%           | 0.187%        |
| 7         | 27.457      | 2684          | 2688        | 2695         | rBV      | 636515         | 1149144       | 8.03%           | 1.445%        |
| 8         | 30.350      | 2999          | 3007        | 3028         | rBV      | 5322063        | 14268207      | 99.73%          | 17.947%       |
| 9         | 30.640      | 3034          | 3039        | 3056         | rVB      | 643855         | 1385965       | 9.69%           | 1.743%        |
| 10        | 34.232      | 3426          | 3435        | 3458         | rBV2     | 4089043        | 11485299      | 80.27%          | 14.446%       |

Sum of corrected areas: 79503938

LSC Report - Integrated Chromatogram

File : C:\MSDCHEM\1\DATA\022702\02577.D  
 Operator : McLean  
 Acquired : 28 Feb 2002 1:29 am using AcqMethod 508SCAN  
 Instrument : Instrumen  
 Sample Name: GC SCAN 2/4 B  
 Misc Info :  
 Vial Number: 20  
 Quant File :5080202.RES (RTE Integrator)



Library Search Compound Report

Data File : C:\MSDCHEM\1\DATA\022702\02577.D  
 Acq On : 28 Feb 2002 1:29 am  
 Sample : GC SCAN 2/4 B  
 Misc :  
 MS Integration Params: LSCINT.P

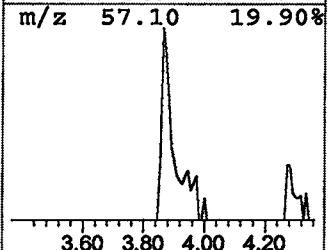
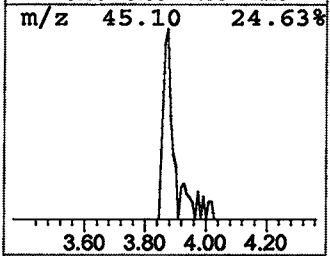
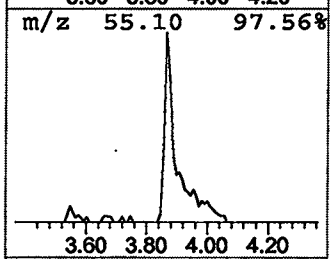
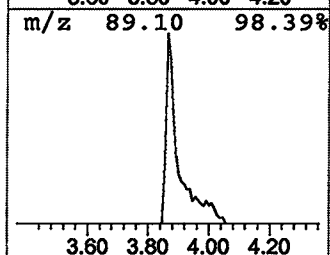
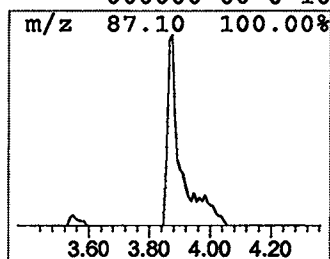
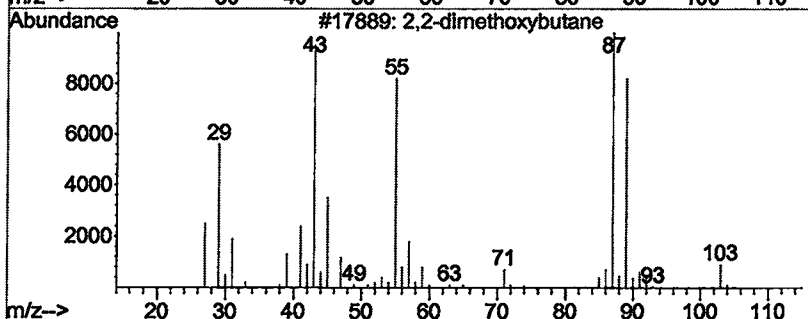
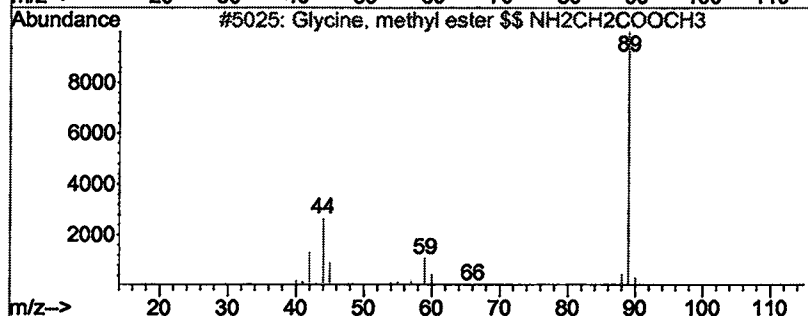
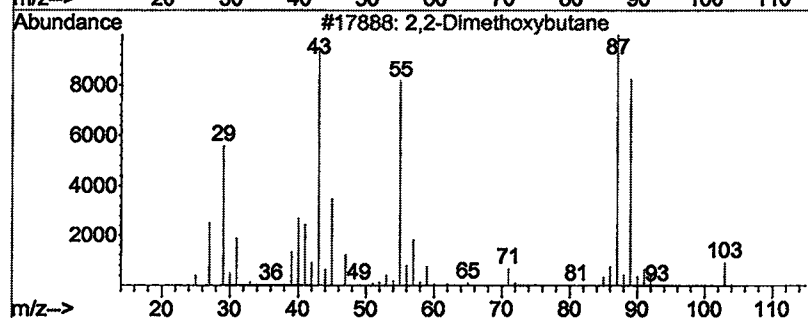
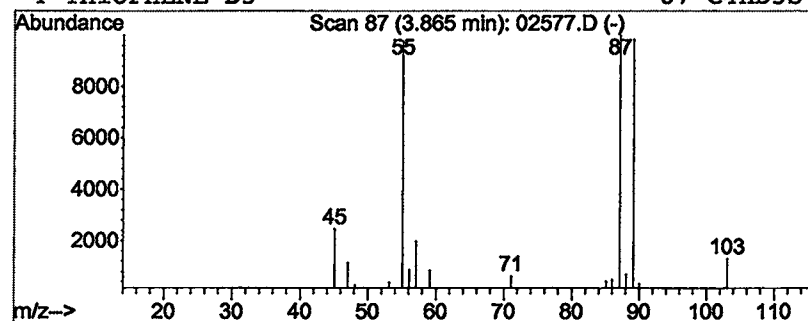
Vial: 20  
 Operator: McLean  
 Inst : Instrumen  
 Multiplr: 1.00

Quant Method : C:\MSDCHEM\1\5973N\5080202.M (RTE Integrator)  
 Title : Quantitation For Method 508gcscan  
 Library : C:\DATABASE\WILEY7N.L

\*\*\*\*\*  
 Peak Number 1 2,2-Dimethoxybutane Concentration Rank 2

| R.T. | EstConc   | Area   | Relative to ISTD       | R.T. |
|------|-----------|--------|------------------------|------|
| 3.87 | 0.42 ug/L | 202285 | 1,4-Dichlorobenzene-d4 | 9.58 |

| Hit# | of | 5 | Tentative ID                          | MW  | MolForm | CAS#        | Qual |
|------|----|---|---------------------------------------|-----|---------|-------------|------|
| 1    |    |   | 2,2-Dimethoxybutane                   | 118 | C6H14O2 | 003453-99-4 | 74   |
| 2    |    |   | Glycine, methyl ester \$\$ NH2CH2C... | 89  | C3H7NO2 | 000616-34-2 | 42   |
| 3    |    |   | 2,2-dimethoxybutane                   | 118 | C6H14O2 | 003453-99-4 | 22   |
| 4    |    |   | THIOPHENE-D3                          | 87  | C4HD3S  | 000000-00-0 | 10   |



## Library Search Compound Report

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Acq On : 28 Feb 2002 1:29 am

Sample : GC SCAN 2/4 B

Misc :

MS Integration Params: LSCINT.P

Vial: 20

Operator: McLean

Inst : Instrumen

Multiplr: 1.00

Quant Method : C:\MSDCHEM\1\5973N\5080202.M (RTE Integrator)

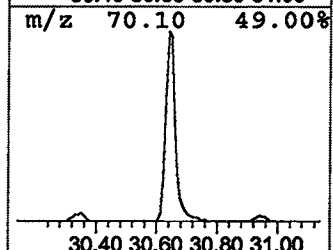
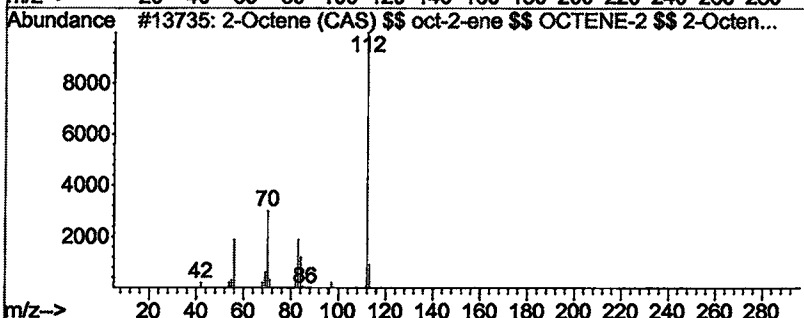
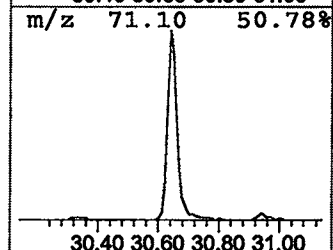
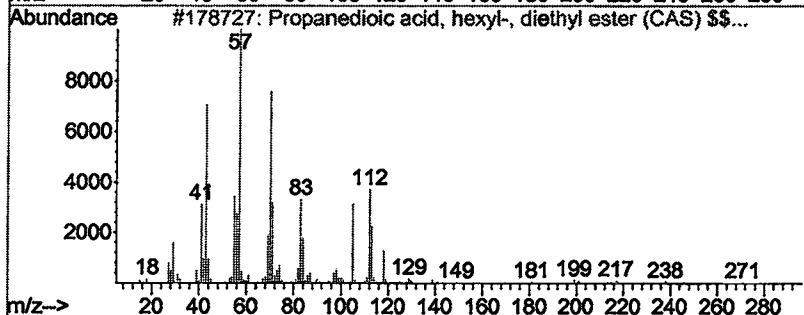
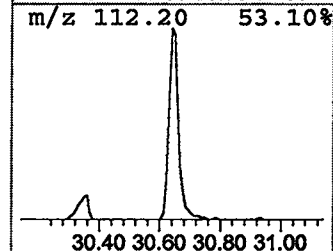
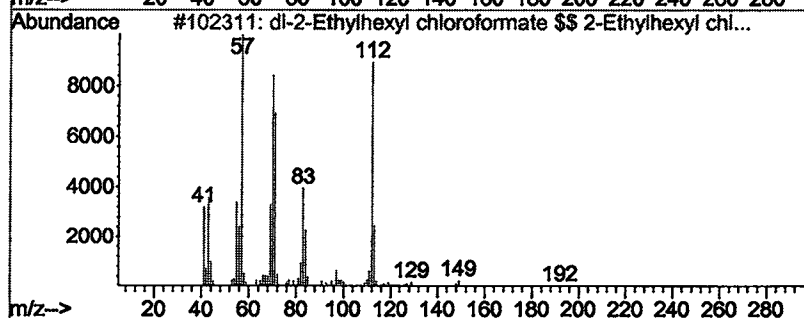
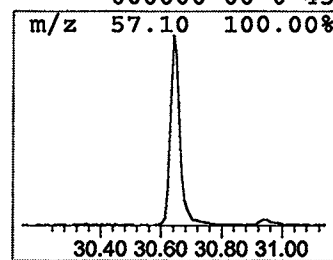
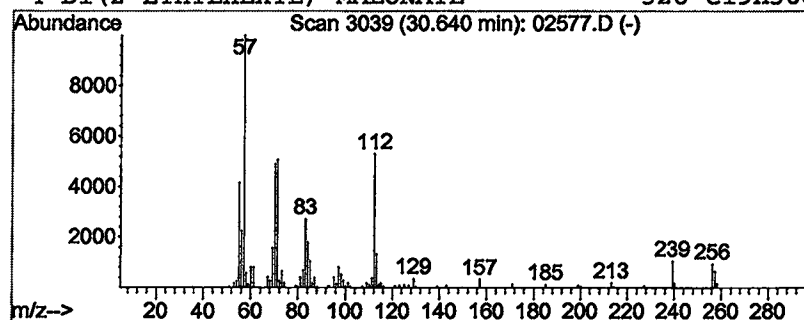
Title : Quantitation For Method 508gcscan

Library : C:\DATABASE\WILEY7N.L

\*\*\*\*\*  
Peak Number 2 dl-2-Ethylhexyl chloroforma... Concentration Rank 1

| R.T.  | EstConc   | Area    | Relative to ISTD | R.T.  |
|-------|-----------|---------|------------------|-------|
| 30.64 | 1.94 ug/L | 1385970 | Chrysene-d12     | 30.35 |

| Hit# | of | 5 | Tentative ID                            | MW  | MolForm   | CAS#        | Qual |
|------|----|---|-----------------------------------------|-----|-----------|-------------|------|
| 1    |    |   | dl-2-Ethylhexyl chloroformate \$\$...   | 192 | C9H17ClO2 | 024468-13-1 | 64   |
| 2    |    |   | Propanedioic acid, hexyl-, dieth...     | 244 | C13H24O4  | 005398-10-7 | 50   |
| 3    |    |   | 2-Octene (CAS) \$\$ oct-2-ene \$\$ O... | 112 | C8H16     | 000111-67-1 | 46   |
| 4    |    |   | DI(2-ETHYLHEXYL) MALONATE               | 328 | C19H36O4  | 000000-00-0 | 43   |



Tentatively Identified Compound (LSC) summary

Operator ID: McLean Date Acquired: 28 Feb 2002 1:29 am  
 Data File: C:\MSDCHEM\1\DATA\022702\02577.D  
 Name: GC SCAN 2/4 B  
 Misc:  
 Method: C:\MSDCHEM\1\5973N\5080202.M (RTE Integrator)  
 Title: Quantitation For Method 508gcscan  
 Library Searched: C:\DATABASE\WILEY7N.L

| TIC Top Hit name     | RT    | EstConc  | Units | Response | --Internal Standard-- |       |          |      |
|----------------------|-------|----------|-------|----------|-----------------------|-------|----------|------|
|                      |       |          |       |          | #                     | RT    | Resp     | Conc |
| 2,2-Dimethoxybutane  | 3.87  | 0.4 ug/L |       | 202285   | 1                     | 9.58  | 9729350  | 20.0 |
| dl-2-Ethylhexyl c... | 30.64 | 1.9 ug/L |       | 1385970  | 5                     | 30.35 | 14268200 | 20.0 |