

Data File : C:\MSDCHEM\1\DATA\022802\01335.D

Acq On : 1 Mar 2002 6:36 am

Sample :

Misc :

MS Integration Params: BENZO.P

Quant Time: Mar 13 06:43:13 2002

Vial: 23

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

*Cherry Mc  
Not needed*

| Internal Standards        | R.T.  | QIon | Response | Conc  | Units | Dev(Min) |
|---------------------------|-------|------|----------|-------|-------|----------|
| 1) 1,4-Dichlorobenzene-d4 | 9.58  | 150  | 1129522  | 20.00 | ug/L  | 0.00     |
| 10) Naphthalene-d8        | 13.05 | 136  | 2965041  | 20.00 | ug/L  | 0.00     |
| 17) Acenaphthene-d10      | 18.16 | 164  | 1652257  | 20.00 | ug/L  | 0.00     |
| 29) Phenanthrene-d10      | 22.52 | 188  | 2558209  | 20.00 | ug/L  | 0.00     |
| 38) Chrysene-d12          | 30.33 | 240  | 2358694  | 20.00 | ug/L  | -0.03    |
| 44) Perylene-d12          | 34.21 | 264  | 1479035  | 20.00 | ug/L  | -0.02    |

## System Monitoring Compounds

|                  |       |     |          |           |      |      |
|------------------|-------|-----|----------|-----------|------|------|
| 37) 2,4-ddd surr | 27.46 | 235 | 216796   | 5.34      | ug/L | 0.00 |
| Spiked Amount    | 5.000 |     | Recovery | = 106.80% |      |      |

## 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         | 18.16 | 165 | 210488 | 9.16 ug/L # | 25     |
| 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        | 0.00  | 149 | 0      | N.D.        |        |
| 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 | 30.94 | 149 | 88463  | 0.83 ug/L   | 96     |
| 43) Chrysene                   | 0.00  | 228 | 0      | N.D. d      |        |
| 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

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Data File : C:\MSDCHEM\1\DATA\022802\01335.D

Vial: 23

Acq On : 1 Mar 2002 6:36 am

Operator: McLean

Sample :

Inst : Instrumen

Misc :

Multiplr: 1.00

MS Integration Params: BENZO.P

Quant Time: Mar 13 06:43:13 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

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Data File : C:\MSDCHEM\1\DATA\022802\01335.D

Acq On : 1 Mar 2002 6:36 am

Sample :

Misc :

MS Integration Params: BENZO.P

Quant Time: Mar 13 6:43 2002

Vial: 23

Operator: McLean

Inst : Instrumen

Multiplr: 1.00

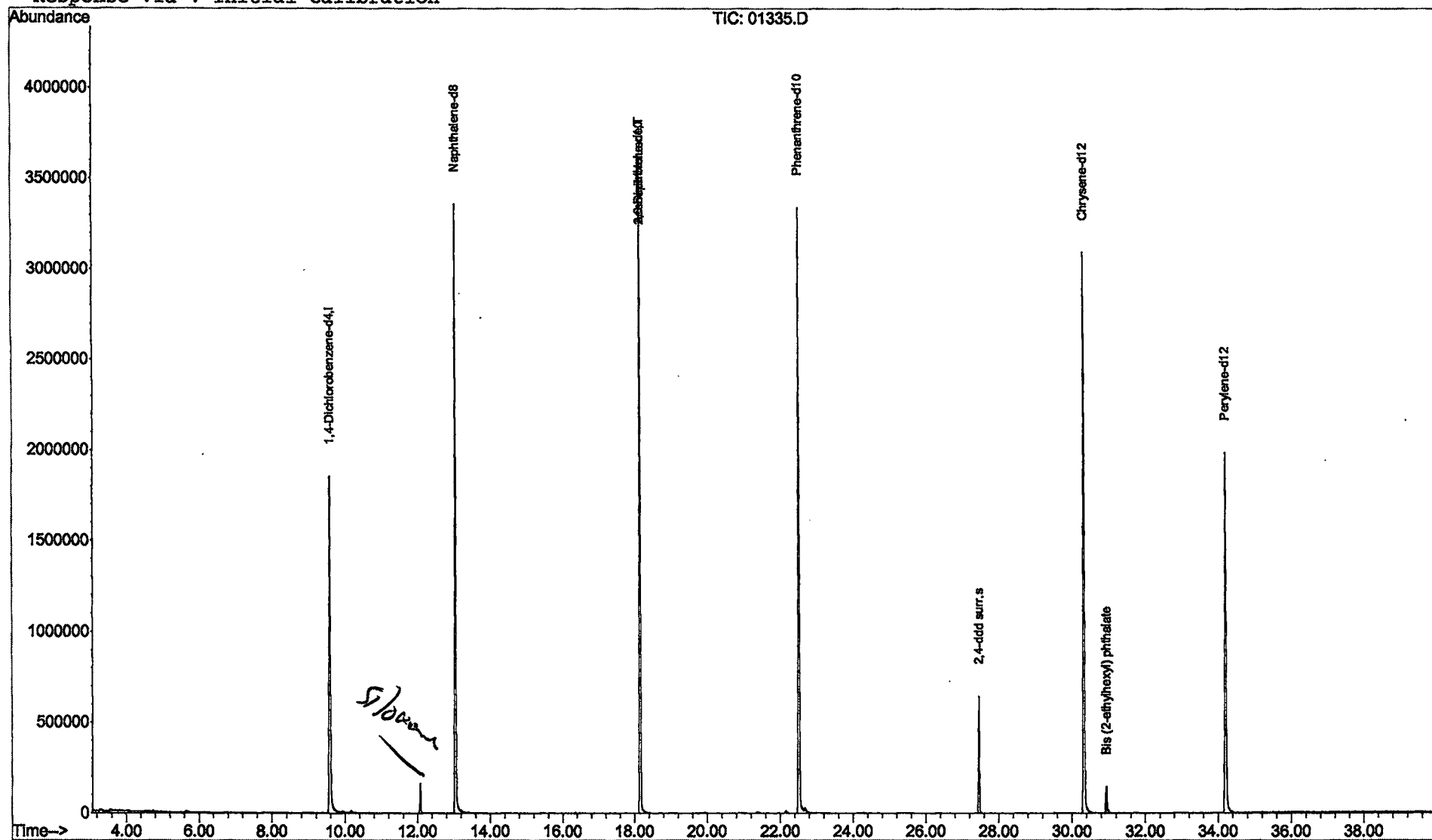
Quant Results File: 5080202.RES

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

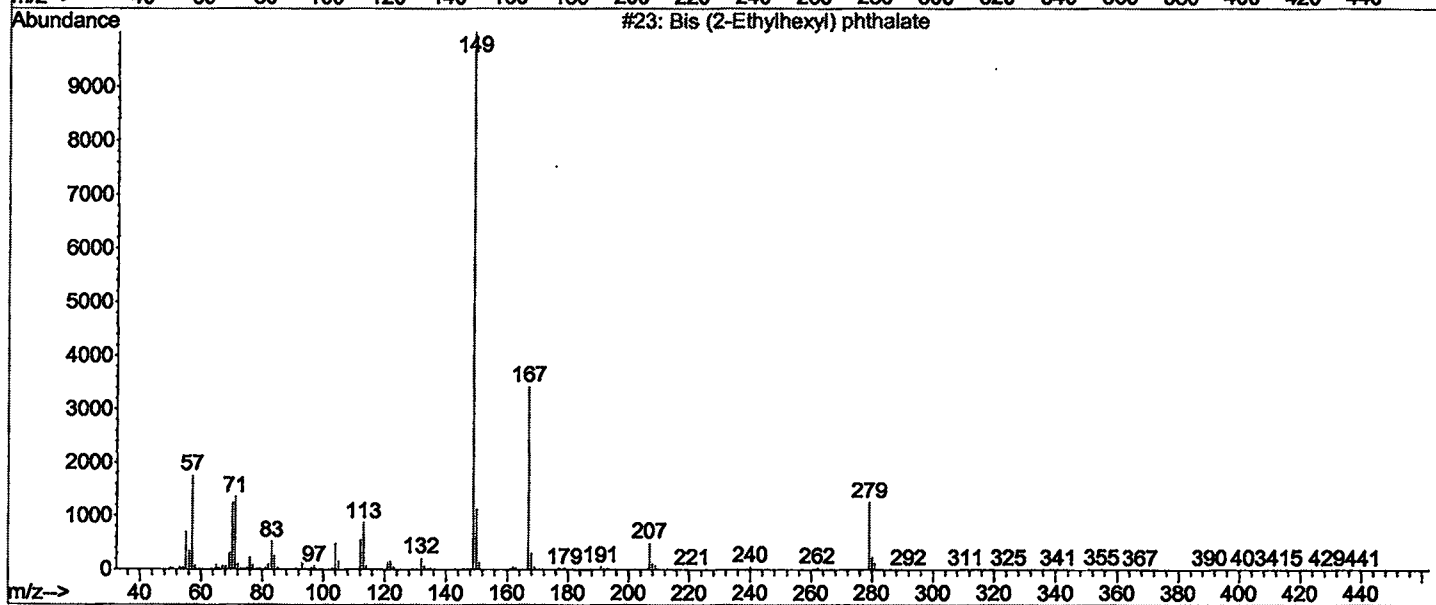
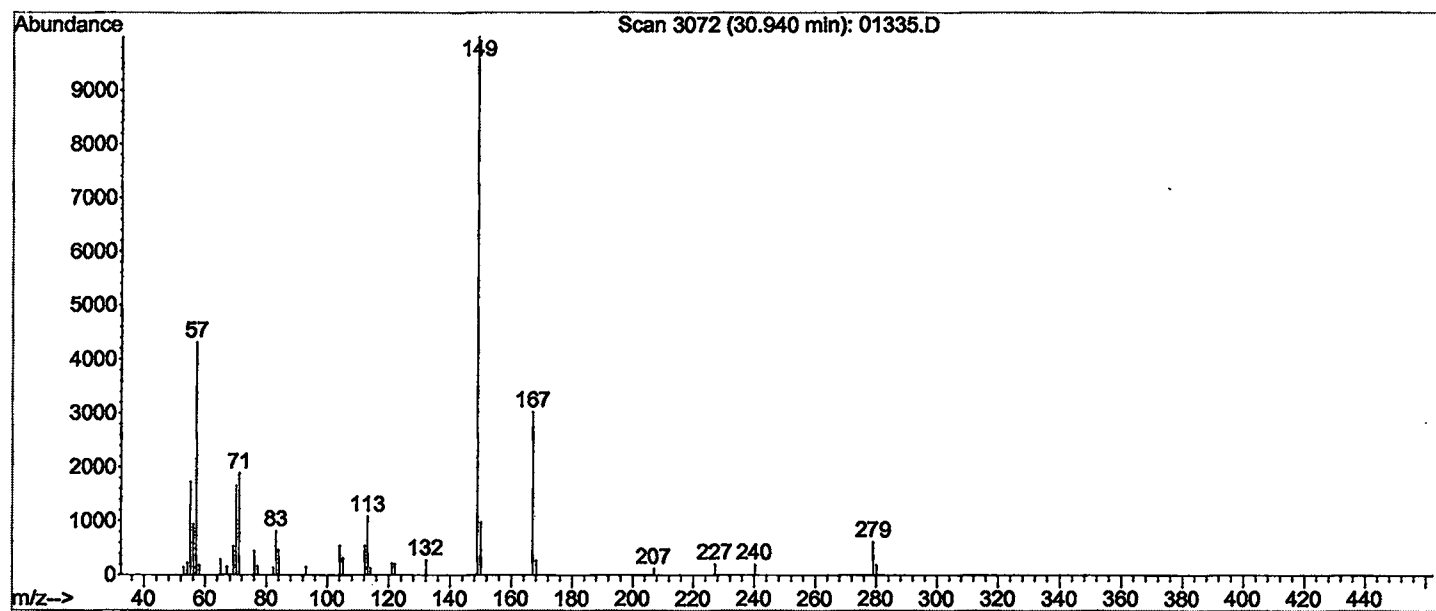


01335.D 5080202.M

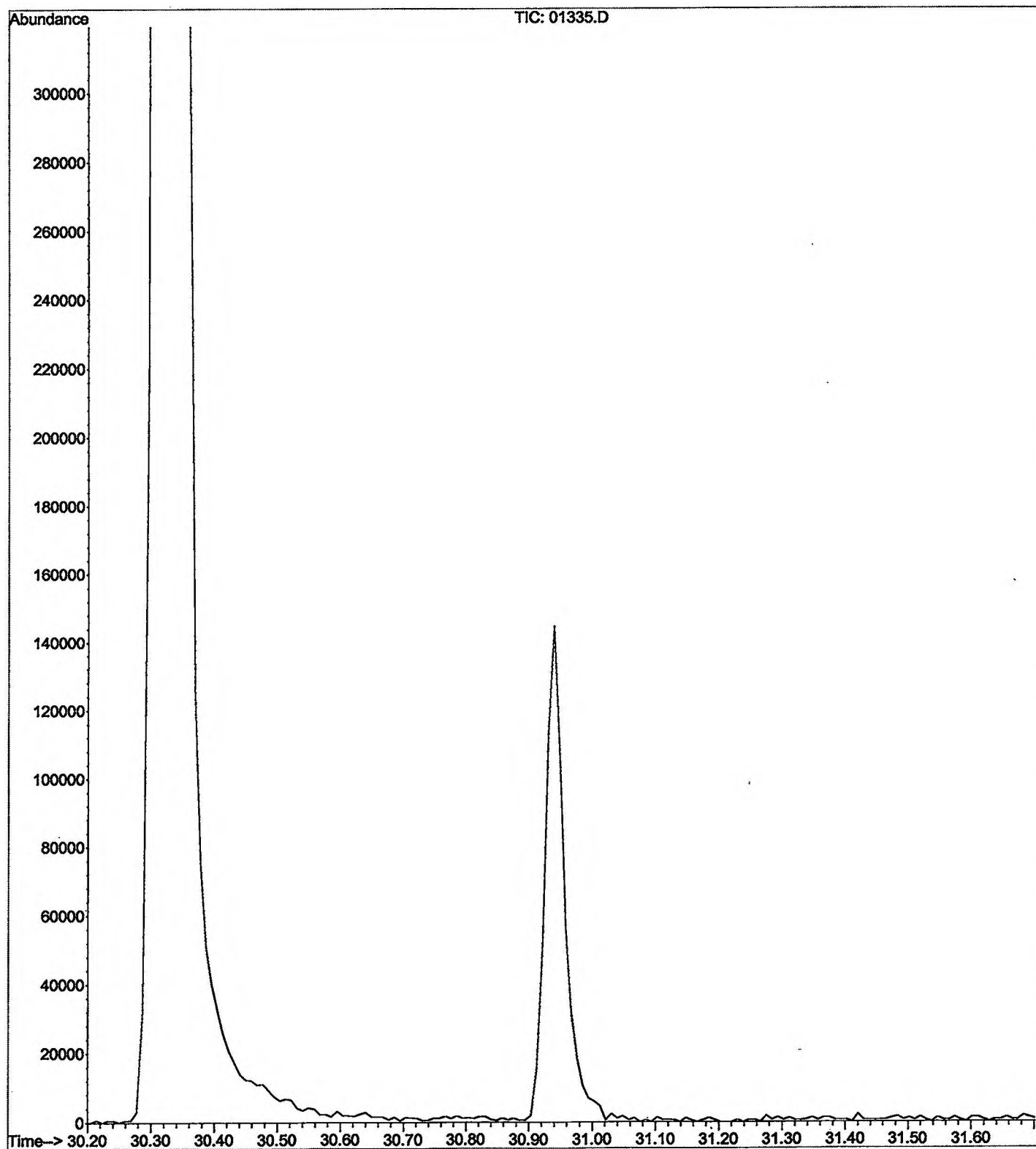
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Library Searched : C:\Database\525.L  
 Quality : 86  
 ID : Bis (2-Ethylhexyl) phthalate



File : C:\MSDCHEM\1\DATA\022802\01335.D  
 Operator : McLean  
 Acquired : 1 Mar 2002 6:36 am using AcqMethod 508SCAN  
 Instrument : Instrumen  
 Sample Name:  
 Misc Info :  
 Vial Number: 23



Data File : C:\MSDCHEM\1\DATA\022802\01335.D

Acq On : 1 Mar 2002 6:36 am

Sample :

Misc :

MS Integration Params: LSCINT.P

Vial: 23

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 &lt; 100 prefer &lt; Baseline drop else tangent &gt;

Peak separation: 5

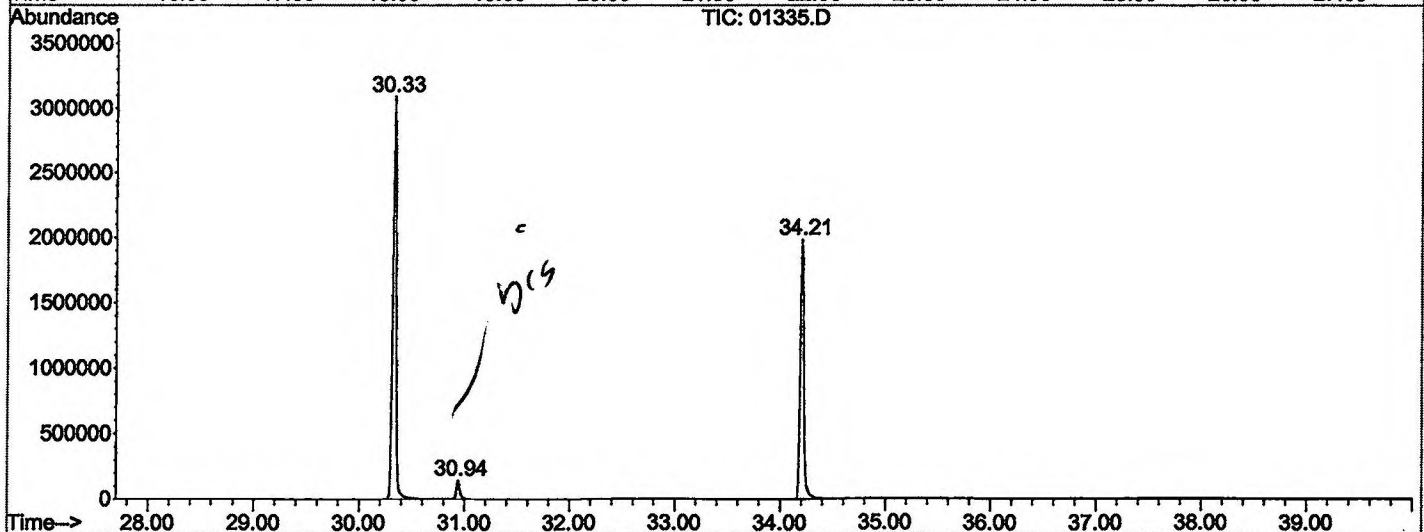
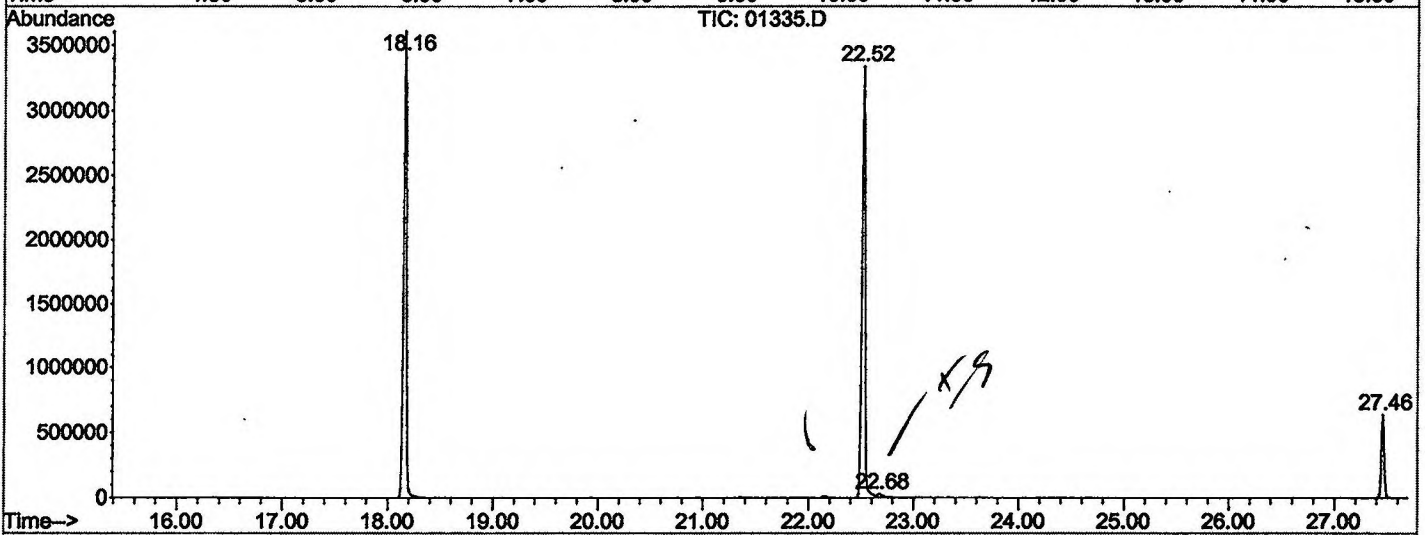
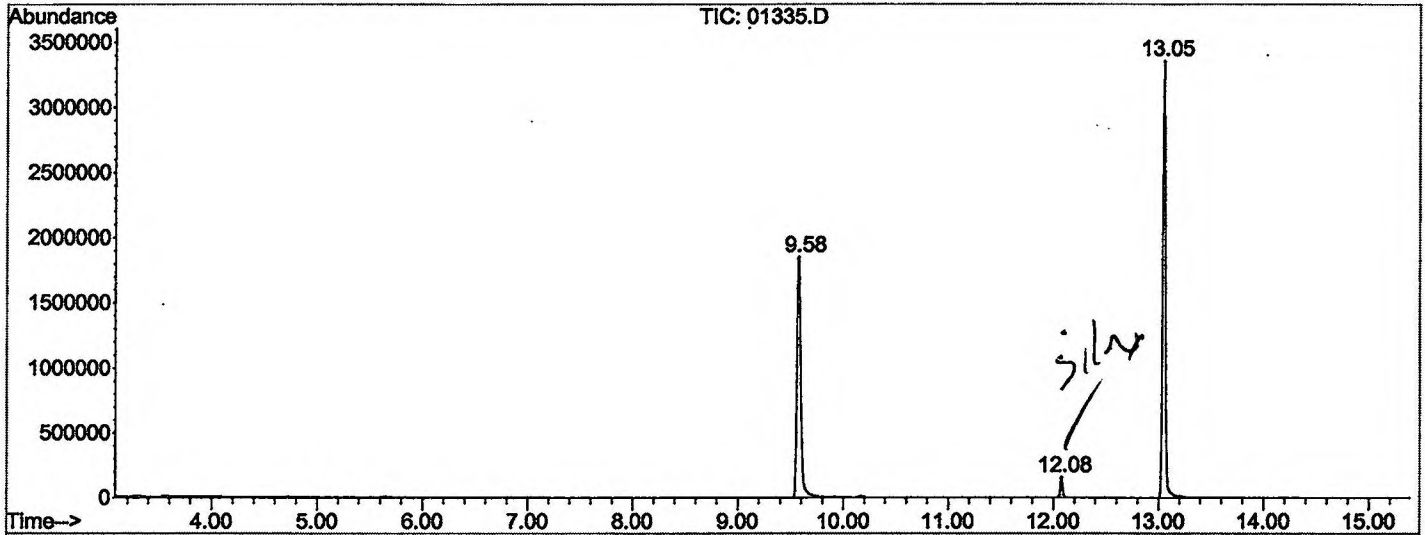
Signal : TIC

| 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         | 9.579       | 712           | 717         | 739          | rBV      | 1854841        | 4755079       | 67.93%          | 12.352%       |
| 2         | 12.083      | 987           | 993         | 999          | rVB2     | 162433         | 268097        | 3.83%           | 0.696%        |
| 3         | 13.053      | 1095          | 1100        | 1114         | rBV      | 3359572        | 6414489       | 91.64%          | 16.663%       |
| 4         | 18.160      | 1657          | 1663        | 1676         | rBV      | 3610643        | 6866668       | 98.10%          | 17.837%       |
| 5         | 22.523      | 2138          | 2144        | 2157         | rBV2     | 3339930        | 6999910       | 100.00%         | 18.184%       |
| 6         | 22.677      | 2157          | 2161        | 2169         | rVB2     | 24124          | 76497         | 1.09%           | 0.199%        |
| 7         | 27.457      | 2683          | 2688        | 2698         | rVB      | 642409         | 1155102       | 16.50%          | 3.001%        |
| 8         | 30.332      | 2998          | 3005        | 3024         | rBV2     | 3092659        | 6802329       | 97.18%          | 17.670%       |
| 9         | 30.940      | 3067          | 3072        | 3081         | rVB      | 144112         | 304330        | 4.35%           | 0.791%        |
| 10        | 34.205      | 3425          | 3432        | 3446         | rBV2     | 1986707        | 4853299       | 69.33%          | 12.607%       |

Sum of corrected areas: 38495800

LSC Report - Integrated Chromatogram

File : C:\MSDCHEM\1\DATA\022802\01335.D  
 Operator : McLean  
 Acquired : 1 Mar 2002 6:36 am using AcqMethod 508SCAN  
 Instrument : Instrumen  
 Sample Name:  
 Misc Info :  
 Vial Number: 23  
 Quant File :5080202.RES (RTE Integrator)



Operator ID: McLean      Date Acquired: 1 Mar 2002      6:36 am  
 Data File: C:\MSDCHEM\1\DATA\022802\01335.D  
 Name:  
 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 |
| Cyclopentasiloxan... | 12.08 | 0.8 ug/L |       | 268097   | 2                     | 13.05 | 6414490 | 20.0 |