

User: Vinci, David L  
Westchester Dept. of Labs & Research  
Date: 04/11/2002 Time: 09:29:48

Sample: AE06053  
Analysis: \$GCMS GCMS Scan For Unknowns  
Units: ug  
Started: 4/11/02 09:29 Ended: 4/11/02 09:29 Analyst: DLV  
Page: 1

|   | Component Name          | Viol | Result | Qualifier | MDL |
|---|-------------------------|------|--------|-----------|-----|
| 1 | Other unknown compounds |      | .      |           |     |
| 2 | Surr_2,4-DDD            |      | 92     |           | 5.0 |

**Comments for Sample AE06053 - Analysis \$GCMS**

Westchester Dept. of Labs & Research

User: Vinci, David L

Date: 04-11-2002 Time: 09:30:04

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\040502\6053.D

Vial: 12

Acq On : 6 Apr 2002 5:26 am

Operator: Bohlk

Sample : SAMPLE 6053 3/15/02

Inst : Instrumen

Misc :

Multiplr: 1.00

MS Integration Params: BENZO.P

Quant Time: Apr 08 13:08:00 2002

Quant Results File: 5080402.RES

Quant Method : C:\MSDCHEM\1\5973N\5080402.M (RTE Integrator)

Title : Quantitation For Method 508gcscan

Last Update : Mon Apr 08 06:47:22 2002

Response via : Initial Calibration

DataAcq Meth : 508SCAN

| Internal Standards        | R.T.  | QIon | Response | Conc  | Units | Dev(Min) |
|---------------------------|-------|------|----------|-------|-------|----------|
| 1) 1,4-Dichlorobenzene-d4 | 9.46  | 152  | 1360387  | 20.00 | ug/L  | -0.01    |
| 10) Naphthalene-d8        | 12.94 | 136  | 6115526  | 20.00 | ug/L  | 0.00     |
| 17) Acenaphthene-d10      | 18.06 | 164  | 3503081  | 20.00 | ug/L  | 0.00     |
| 29) Phenanthrene-d10      | 22.42 | 188  | 4925533  | 20.00 | ug/L  | 0.00     |
| 38) Chrysene-d12          | 30.24 | 240  | 3872077  | 20.00 | ug/L  | -0.01    |
| 44) Perylene-d12          | 34.12 | 264  | 2451701  | 20.00 | ug/L  | 0.00     |

## System Monitoring Compounds

|                    |       |     |          |      |        |      |
|--------------------|-------|-----|----------|------|--------|------|
| 37) 2,4-DDD (SURR) | 27.36 | 235 | 225384   | 4.61 | ug/L   | 0.00 |
| Spiked Amount      | 5.000 |     | Recovery | =    | 92.20% |      |

## 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. | d       |
| 34) Anthracene                 | 0.00  | 178 | 0    | N.D. | d       |
| 35) Di-n-butylphthalate        | 0.00  | 149 | 0    | N.D. | 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 | 30.85 | 149 | 7877 | 0.06 | ug/L 90 |
| 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

6053.D 5080402.M Mon Apr 08 13:09:54 2002

Page 1

## Quantitation Report (QT Reviewed)

Data File : C:\MSDCHEM\1\DATA\040502\6053.D

Vial: 12

Acq On : 6 Apr 2002 5:26 am

Operator: Bohlk

Sample : SAMPLE 6053 3/15/02

Inst : Instrumen

Misc :

Multiplr: 1.00

MS Integration Params: BENZO.P

Quant Time: Apr 08 13:08:00 2002

Quant Results File: 5080402.RES

Quant Method : C:\MSDCHEM\1\5973N\5080402.M (RTE Integrator)

Title : Quantitation For Method 508gcscan

Last Update : Mon Apr 08 06:47:22 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

6053.D 5080402.M Mon Apr 08 13:09:54 2002

Data File : C:\MSDCHEM\1\DATA\040502\6053.D

Vial: 12

Acq On : 6 Apr 2002 5:26 am

Operator: Bohlk

Sample : SAMPLE 6053 3/15/02

Inst : Instrumen

Misc :

Multiplr: 1.00

MS Integration Params: BENZO.P

Quant Time: Apr 8 13:09 2002

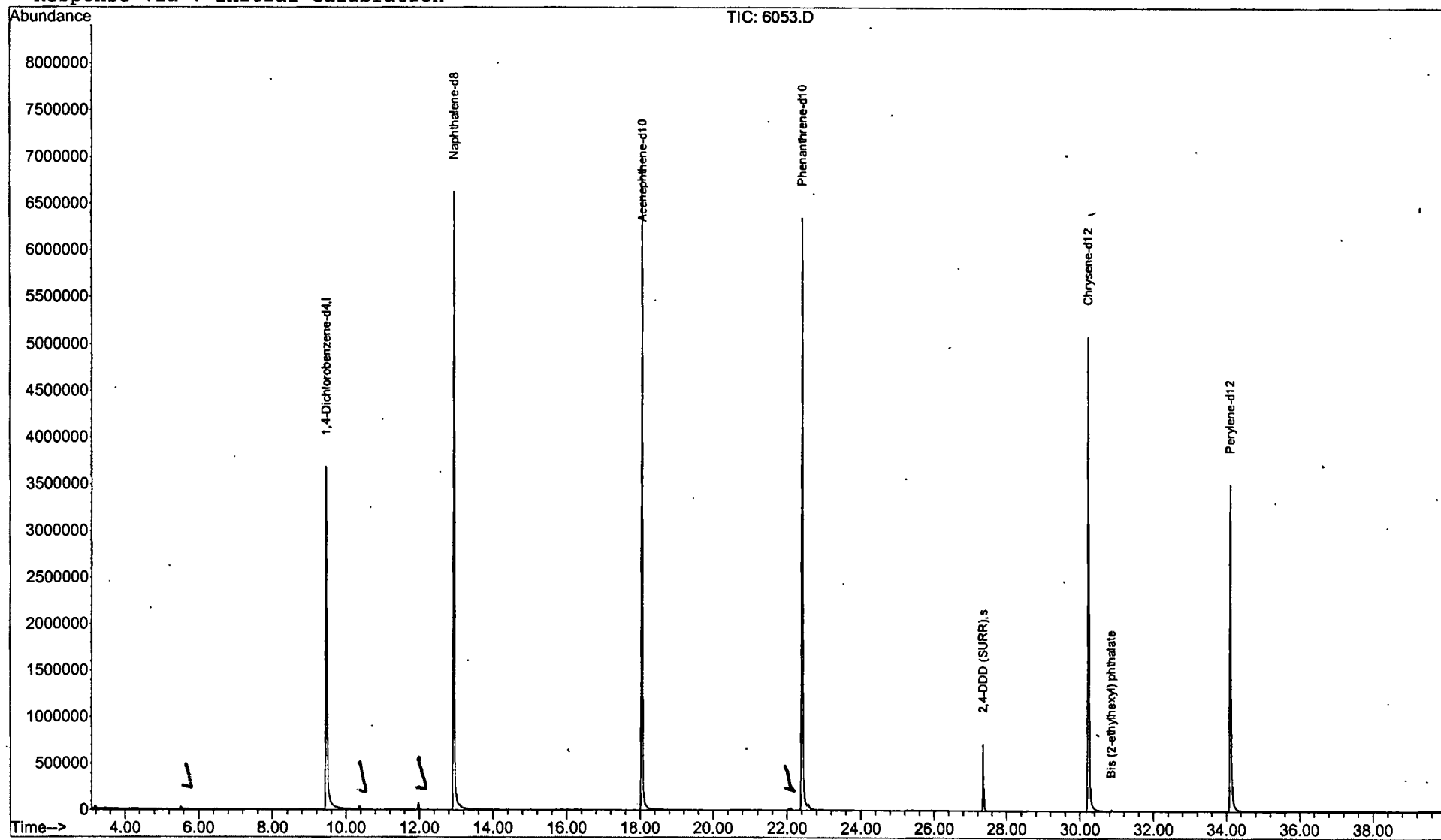
Quant Results File: 5080402.RES

Method : C:\MSDCHEM\1\5973N\5080402.M (RTE Integrator)

Title : Quantitation For Method 508gcscan

Last Update : Mon Apr 08 06:47:22 2002

Response via : Initial Calibration



6053.D 5080402.M

Mon Apr 08 13:09:55 2002

Page 3

## LSC Area Percent Report

Data File : C:\MSDCHEM\1\DATA\040502\6053.D

Acq On : 6 Apr 2002 5:26 am

Sample : SAMPLE 6053 3/15/02

Misc :

MS Integration Params: LSCINT.P

Vial: 12

Operator: Bohlk

Inst : Instrumen

Multiplr: 1.00

Method : C:\MSDCHEM\1\5973N\5080402.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

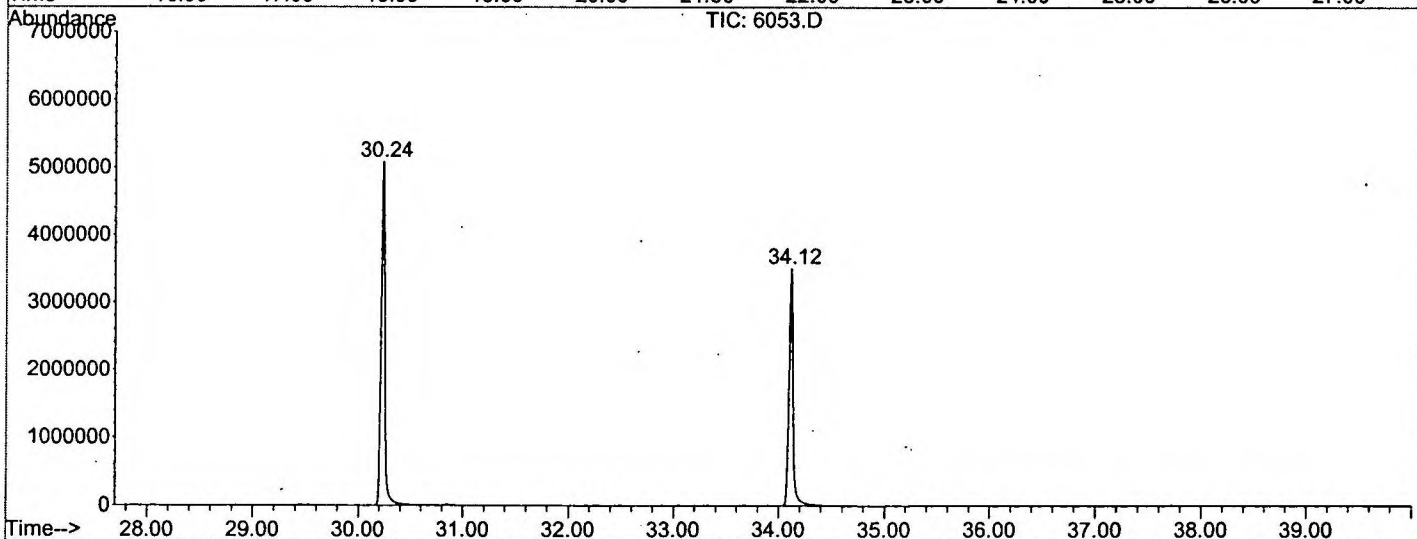
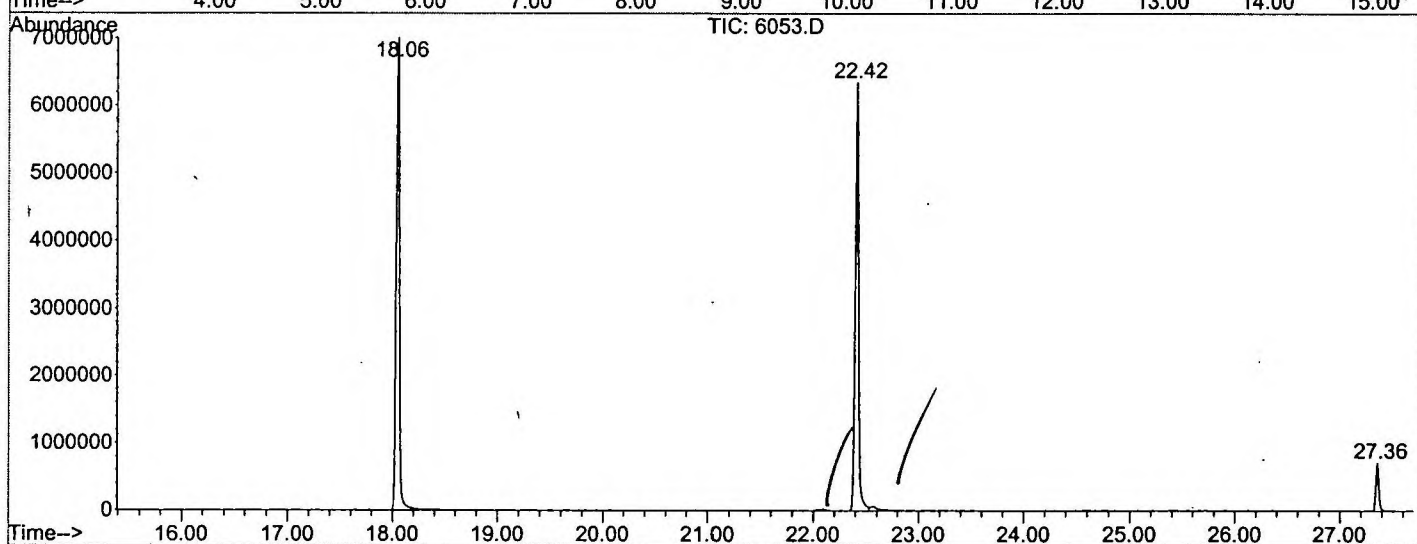
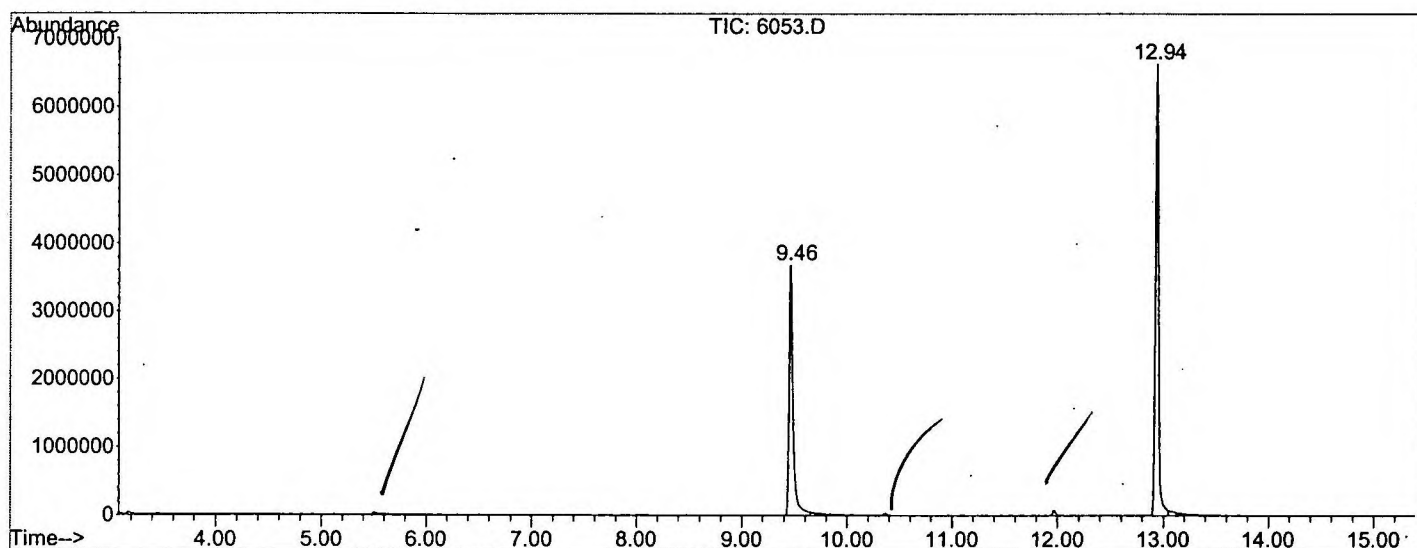
Signal : TIC

| 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         | 9.460       | 1079          | 1086        | 1118         | rBV      | 3676599        | 9337271       | 61.64%          | 12.383%       |
| 2         | 12.938      | 1669          | 1678        | 1696         | rBV      | 6624968        | 13465234      | 88.89%          | 17.857%       |
| 3         | 18.056      | 2539          | 2549        | 2577         | rBV      | 6999380        | 15148523      | 100.00%         | 20.089%       |
| 4         | 22.416      | 3280          | 3291        | 3310         | rBV2     | 6347305        | 14393473      | 95.02%          | 19.088%       |
| 5         | 27.357      | 4124          | 4132        | 4147         | rBV2     | 711546         | 1354487       | 8.94%           | 1.796%        |
| 6         | 30.236      | 4610          | 4622        | 4645         | rBV2     | 5075461        | 12427792      | 82.04%          | 16.481%       |
| 7         | 34.120      | 5271          | 5283        | 5308         | rBV2     | 3502550        | 9279868       | 61.26%          | 12.306%       |

Sum of corrected areas: 75406648

# LSC Report - Integrated Chromatogram

File : C:\MSDCHEM\1\DATA\040502\6053.D  
 Operator : Bohlk  
 Acquired : 6 Apr 2002 5:26 am using AcqMethod 508SCAN  
 Instrument : Instrumen  
 Sample Name: SAMPLE 6053 3/15/02  
 Misc Info :  
 Vial Number: 12  
 Quant File :5080402.RES (RTE Integrator)



Tentatively Identified Compound (LSC) summary

Operator ID: Bohlk      Date Acquired: 6 Apr 2002    5:26 am  
 Data File: C:\MSDCHEM\1\DATA\040502\6053.D  
 Name: SAMPLE 6053    3/15/02  
 Misc:  
 Method: C:\MSDCHEM\1\5973N\5080402.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 |
| -----            |    |         |       |          |                       |    |      |      |