

User: Vinci, David L  
 Westchester Dept. of Labs & Research  
 Date: 02/25/2002 Time: 15:55:25

Sample: AE00109  
 Analysis: \$GCMS GCMS Scan For Unknowns  
 Units: ug  
 Started: 2/25/02 15:54 Ended: 2/25/02 15:54 Analyst: DLV  
 Page: 1

|   | Component Name          | Viol | Result | Qualifier | MDL |
|---|-------------------------|------|--------|-----------|-----|
| 1 | Other unknown compounds |      | .      |           |     |
| 2 | Surr_2,4-DDD            |      | 79     |           | 5.0 |

**Comments for Sample AE00109 - Analysis \$GCMS**

Westchester Dept. of Labs & Research

User: Vinci, David L

Date: 02-25-2002 Time: 15:56:45

The GCMS scan, from 35 to 550 amu, does not reveal the presence of unusual compounds. The recoveries for 11 of the 43 compounds in the LCS Standard were below their acceptance ranges.

## Quantitation Report

(QT Reviewed)

Data File : C:\HPCHEM\1\DATA\FEB1402\109.D  
 Acq On : 15 Feb 2002 9:18 am  
 Sample : GC/MS SCAN 2/11  
 Misc :  
 MS Integration Params: GOOFY.P  
 Quant Time: Feb 15 12:09 2002

Vial: 32  
 Operator: 209 GALOS  
 Inst : GC/MS 209  
 Multiplr: 1.00

Quant Results File: GCMS0208.RES

Quant Method : C:\HPCHEM\1\METHODS\GCMS0208.M (RTE Integrator)  
 Title : 209 625/TCLP calibration table  
 Last Update : Wed Feb 13 11:43:23 2002  
 Response via : Initial Calibration  
 DataAcq Meth : GCMS0208

| Internal Standards        | R.T.  | QIon | Response | Conc  | Units | Dev (Min) |
|---------------------------|-------|------|----------|-------|-------|-----------|
| 1) 1,4-Dichlorobenzene-d4 | 12.35 | 152  | 359382   | 20.00 | ug/l  | 0.00      |
| 10) Naphthalene-d8        | 15.99 | 136  | 1409055  | 20.00 | ug/l  | 0.00      |
| 17) Acenaphthene-d10      | 21.31 | 164  | 759623   | 20.00 | ug/l  | 0.00      |
| 29) Phenanthrene-d10      | 25.75 | 188  | 1129779  | 20.00 | ug/l  | -0.02     |
| 38) Chrysene-d12          | 33.82 | 240  | 821437   | 20.00 | ug/l  | -0.02     |
| 44) Perylene-d12          | 38.11 | 264  | 531214   | 20.00 | ug/l  | -0.02     |

## System Monitoring Compounds

|               |       |     |          |      |        |      |
|---------------|-------|-----|----------|------|--------|------|
| 37) 2,4-DDD   | 30.83 | 235 | 44942    | 3.96 | ug/mL  | 0.33 |
| Spiked Amount | 5.000 |     | Recovery | =    | 79.20% |      |

## Target Compounds

|                                 |       |     |     | Qvalue |
|---------------------------------|-------|-----|-----|--------|
| 2) N-nitroso-dimethyl amine     | 0.00  | 74  | 0   | N.D.   |
| 3) bis(2-Chloroethyl) ether     | 0.00  | 93  | 0   | N.D.   |
| 4) 1,3-Dichlorobenzene          | 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-Chloropropan   | 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) methane | 0.00  | 93  | 0   | N.D.   |
| 14) 1,2,4-Trichlorobenzene      | 0.00  | 180 | 0   | N.D.   |
| 15) Naphthalene                 | 16.05 | 128 | 289 | 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.   |
| 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-phenylether  | 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.   |

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

109.D GCMS0208.M Fri Feb 15 12:10:32 2002

Page 1

## Quantitation Report

(QT Reviewed)

Data File : C:\HPCHEM\1\DATA\FEB1402\109.D

Vial: 32

Acq On : 15 Feb 2002 9:18 am

Operator: 209 GALOS

Sample : GC/MS SCAN 2/11

Inst : GC/MS 209

Misc :

Multiplr: 1.00

MS Integration Params: GOOFY.P

Quant Time: Feb 15 12:09 2002

Quant Results File: GCMS0208.RES

Quant Method : C:\HPCHEM\1\METHODS\GCMS0208.M (RTE Integrator)

Title : 209 625/TCLP calibration table

Last Update : Wed Feb 13 11:43:23 2002

Response via : Initial Calibration

DataAcq Meth : GCMS0208

| Compound                       | R.T.  | QIon | Response | Conc | Unit | Qvalue |
|--------------------------------|-------|------|----------|------|------|--------|
| 30) N-Nitrosodiphenylamine     | 0.00  | 168  | 0        | N.D. |      |        |
| 31) 4-Bromophenyl-phenylether  | 0.00  | 248  | 0        | N.D. |      |        |
| 32) Hexachlorobenzene          | 0.00  | 284  | 0        | N.D. |      |        |
| 33) Phenanthrene               | 25.76 | 178  | 320      | N.D. |      |        |
| 34) Anthracene                 | 25.76 | 178  | 320      | N.D. |      |        |
| 35) Di-n-butylphthalate        | 27.72 | 149  | 10130    | 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         | 33.83 | 228  | 1966     | N.D. |      |        |
| 42) Bis(2-ethylhexyl)phthalate | 34.02 | 149  | 4265     | N.D. |      |        |
| 43) Chrysene                   | 33.83 | 228  | 1966     | N.D. |      |        |
| 45) Di-n-Octylphthalate        | 0.00  | 149  | 0        | N.D. |      |        |
| 46) Benzo(b)fluoranthene       | 0.00  | 252  | 0        | N.D. |      |        |
| 47) Benzo(k)fluoranthene       | 0.00  | 252  | 0        | N.D. |      |        |
| 48) Benzo(a)pyrene             | 38.12 | 252  | 2079     | N.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

109.D GCMS0208.M

Fri Feb 15 12:10:36 2002

Page 2

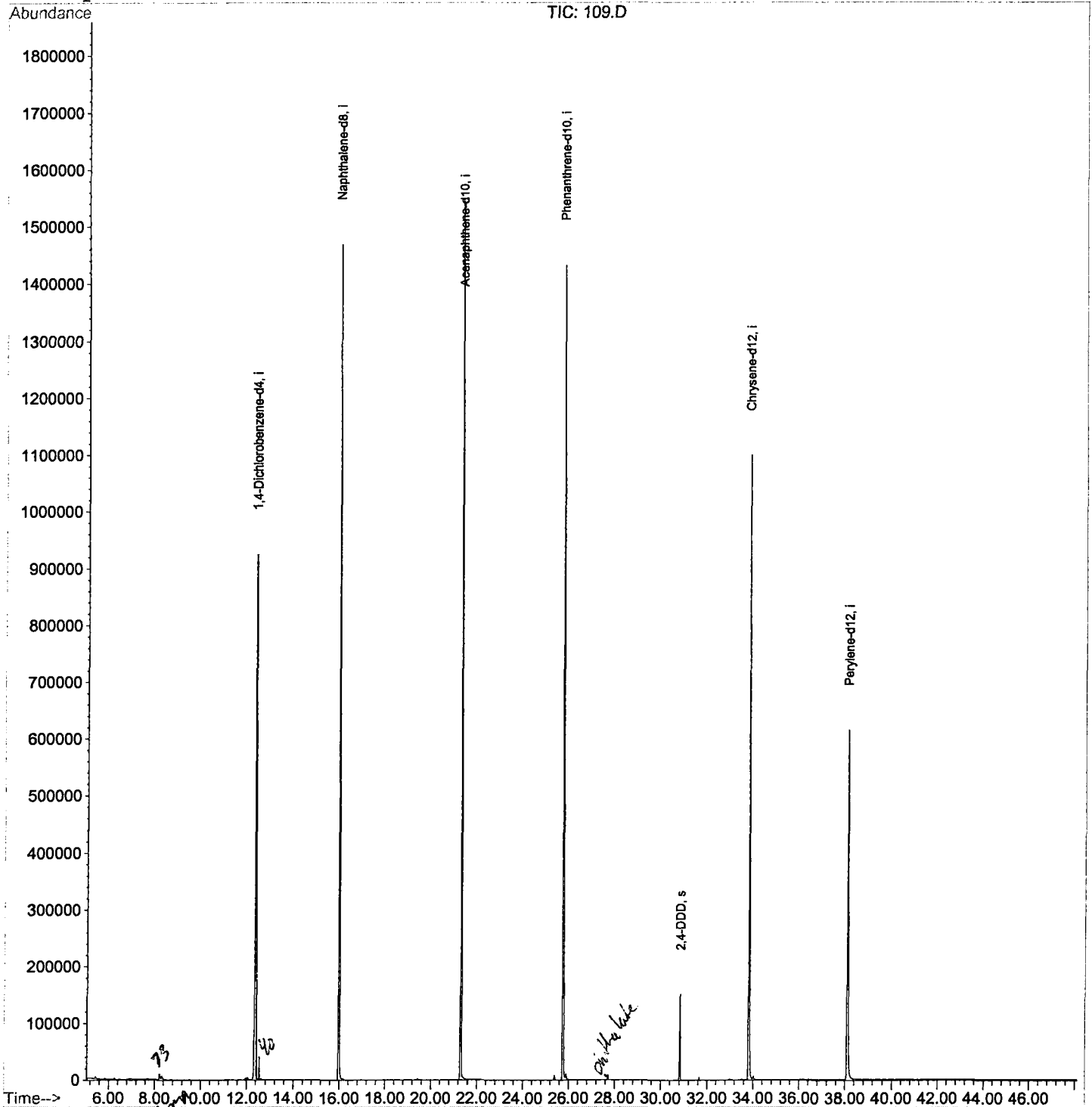
# Quantitation Report

Data File : C:\HPCHEM\1\DATA\FEB1402\109.D  
 Acq On : 15 Feb 2002 9:18 am  
 Sample : GC/MS SCAN 2/11  
 Misc :  
 MS Integration Params: GOOFY.P  
 Quant Time: Feb 15 12:09 2002

Vial: 32  
 Operator: 209 GALOS  
 Inst : GC/MS 209  
 Multiplr: 1.00

Quant Results File: GCMS0208.RES

Method : C:\HPCHEM\1\METHODS\GCMS0208.M (RTE Integrator)  
 Title : 209 625/TCLP calibration table  
 Last Update : Wed Feb 13 11:43:23 2002  
 Response via : Initial Calibration



## LSC Area Percent Report

Data File : C:\HPCHEM\1\DATA\FEB1402\109.D  
 Acq On : 15 Feb 2002 9:18 am  
 Sample : GC/MS SCAN 2/11  
 Misc :  
 MS Integration Params: LSCINT.P

Vial: 32  
 Operator: 209 GALOS  
 Inst : GC/MS 209  
 Multiplr: 1.00

Method : C:\HPCHEM\1\METHODS\GCMS0208.M (RTE Integrator)  
 Title : 209 625/TCLP calibration table  
 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

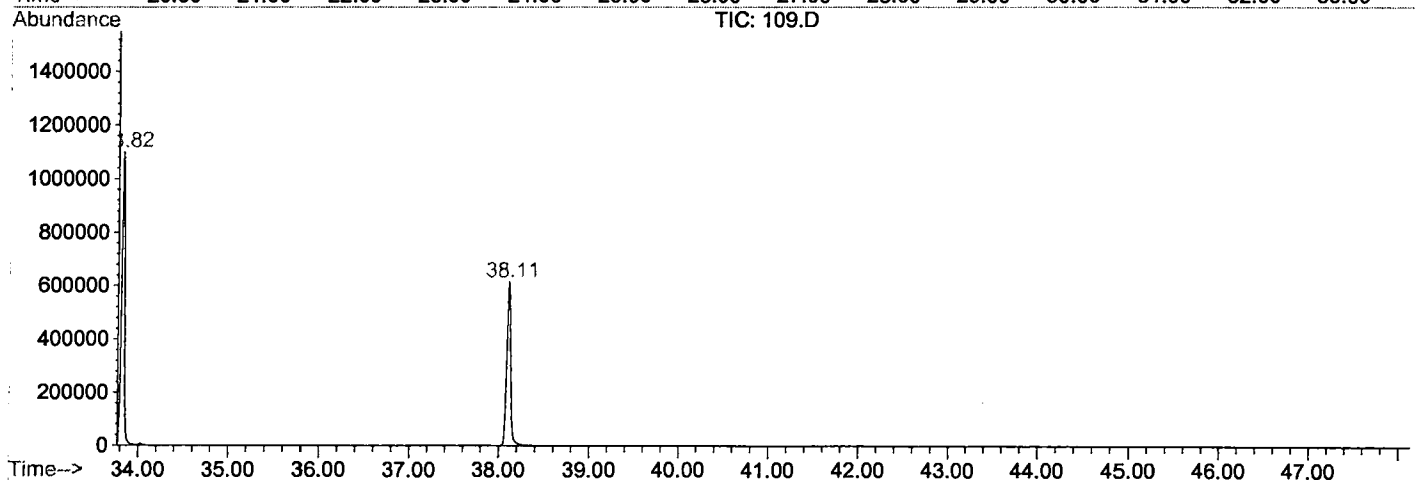
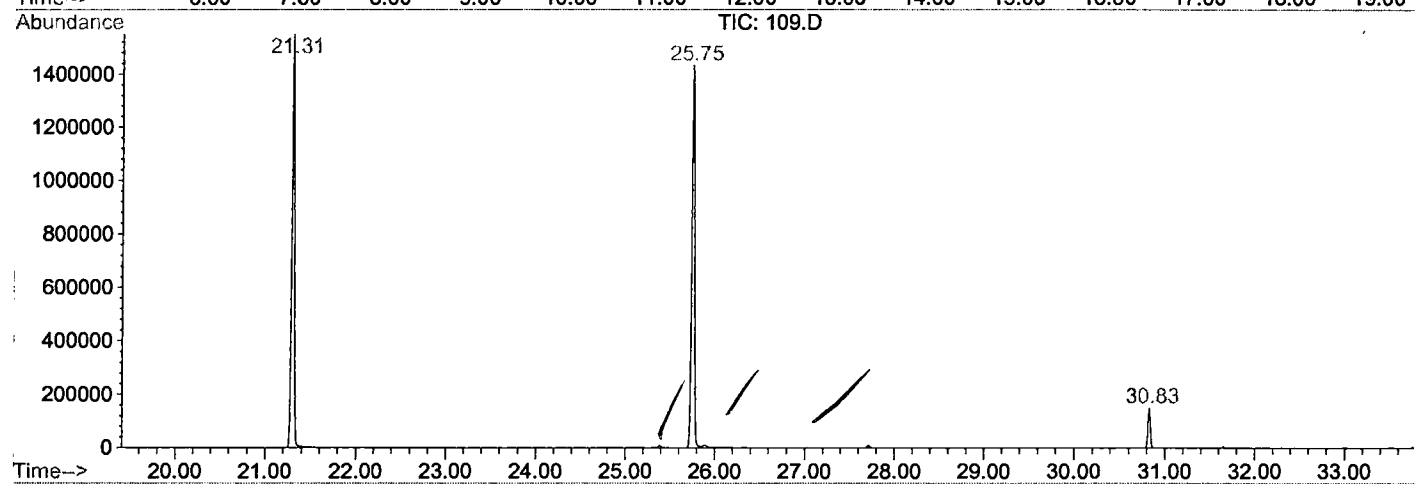
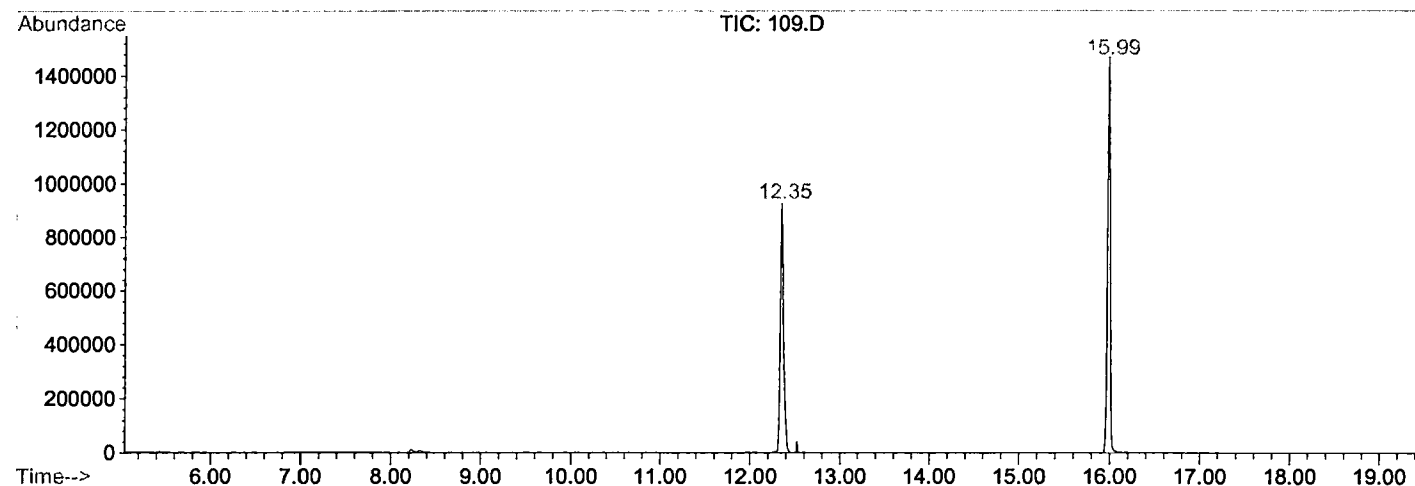
| peak<br># | R.T.<br>min | first<br>scan | max<br>scan | last<br>scan | PK<br>TY | peak<br>height | peak<br>area | peak<br>% max. | % of<br>total |
|-----------|-------------|---------------|-------------|--------------|----------|----------------|--------------|----------------|---------------|
| 1         | 12.348      | 790           | 796         | 810          | rVB      | 924426         | 2342486      | 74.53%         | 14.501%       |
| 2         | 15.992      | 1186          | 1193        | 1210         | rBV      | 1469058        | 2930641      | 93.25%         | 18.142%       |
| 3         | 21.308      | 1765          | 1772        | 1789         | rBV      | 1549036        | 3142834      | 100.00%        | 19.455%       |
| 4         | 25.751      | 2249          | 2256        | 2267         | rBV2     | 1432185        | 3131989      | 99.65%         | 19.388%       |
| 5         | 30.828      | 2804          | 2809        | 2814         | rBV      | 150540         | 274834       | 8.74%          | 1.701%        |
| 6         | 33.820      | 3128          | 3135        | 3153         | rBV2     | 1098004        | 2494773      | 79.38%         | 15.444%       |
| 7         | 38.108      | 3593          | 3602        | 3623         | rBV2     | 613221         | 1836509      | 58.43%         | 11.369%       |

Sum of corrected areas: 16154066

109.D GCMS0208.M Fri Feb 15 13:59:59 2002

# LSC Report - Integrated Chromatogram

File : C:\HPCHEM\1\DATA\FEB1402\109.D  
 Operator : 209 GALOS  
 Acquired : 15 Feb 2002 9:18 am using AcqMethod GCMS0208  
 Instrument : GC/MS 209  
 Sample Name: GC/MS SCAN 2/11  
 Misc Info :  
 Vial Number: 32  
 Quant File :GCMS0208.RES (RTE Integrator)



109.D GCMS0208.M

Fri Feb 15 14:00:04 2002

Tentatively Identified Compound (LSC) summary

Operator ID: 209 GALOS Date Acquired: 15 Feb 2002 9:18 am  
 Data File: C:\HPCHEM\1\DATA\FEB1402\109.D  
 Name: GC/MS SCAN 2/11  
 Misc:  
 Method: C:\HPCHEM\1\METHODS\GCMS0208.M (RTE Integrator)  
 Title: 209 625/TCLP calibration table  
 Library Searched: C:\DATABASE\NBS75K.L

| TIC Top Hit name | RT         | EstConc  | Units | Area | IntStd | ISRT | ISArea | ISConc |
|------------------|------------|----------|-------|------|--------|------|--------|--------|
| -----            |            |          |       |      |        |      |        |        |
| 109.D GCMS0208.M | Fri Feb 15 | 14:00:09 | 2002  |      |        |      |        |        |