

User: Galos, Marie  
 Westchester Dept. of Labs & Research  
 Date: 12/03/2001 Time: 15:00:28

Sample: AD28593  
 Analysis: \$C1GCMS CALI GCMS  
 Units: ug  
 Started: 11/26/20 00:01 Ended: 11/26/20 00:01 Analyst: MG  
 Result source: m:\instruments\209\data\nov2601\cal25.d\cal25.rr  
 Page: 1

|    | Component Name                   | Viol | Result | MDL |
|----|----------------------------------|------|--------|-----|
| 1  | n-Nitrosodimethylamine           |      | 25     | 2.0 |
| 2  | bis(2-Chloroethyl)ether          |      | 24     | 2.0 |
| 3  | 1,3-Dichlorobenzene              |      | 25     | 2.0 |
| 4  | 1,4-Dichlorobenzene              |      | 25     | 2.0 |
| 5  | 1,2-Dichlorobenzene              |      | 25     | 2.0 |
| 6  | 2,2'-oxybis(1-Chloropropane)     |      | 25     | 2.0 |
| 7  | n-Nitrosodi-n-propylamine        |      | 25     | 2.0 |
| 8  | Hexachloroethane                 |      | 25     | 2.0 |
| 9  | Nitrobenzene                     |      | 25     | 2.0 |
| 10 | Isophorone                       |      | 25     | 2.0 |
| 11 | bis(2-Chloroethoxy)methane       |      | 25     | 2.0 |
| 12 | 1,2,4-Trichlorobenzene           |      | 25     | 2.0 |
| 13 | Naphthalene                      |      | 25     | 2.0 |
| 14 | Hexachlorobutadiene              |      | 25     | 2.0 |
| 15 | Hexachlorocyclopentadiene        |      | 23     | 2.0 |
| 16 | 2-Chloronaphthalene              |      | 25     | 2.0 |
| 17 | Dimethylphthalate                |      | 25     | 2.0 |
| 18 | 2,6-Dinitrotoluene               |      | 25     | 2.0 |
| 19 | Acenaphthylene                   |      | 25     | 2.0 |
| 20 | Acenaphthene                     |      | 25     | 2.0 |
| 21 | 2,4-Dinitrotoluene               |      | 25     | 2.0 |
| 22 | Diethylphthalate                 |      | 25     | 2.0 |
| 23 | 4-Chlorophenylphenylether        |      | 25     | 2.0 |
| 24 | Fluorene                         |      | 25     | 2.0 |
| 25 | Azobenzene/1,2-Diphenylhydrazine |      | 24     | 2.0 |
| 26 | n-Nitrosodiphenylamine           |      | 24     | 2.0 |
| 27 | 4-Bromophenylphenylether         |      | 25     | 2.0 |
| 28 | Hexachlorobenzene                |      | 25     | 2.0 |
| 29 | Phenanthrene                     |      | 25     | 2.0 |
| 30 | Anthracene                       |      | 25     | 2.0 |
| 31 | Di-n-butylphthalate              |      | 25     | 2.0 |
| 32 | Fluoranthene                     |      | 25     | 2.0 |
| 33 | Pyrene                           |      | 25     | 2.0 |
| 34 | Butylbenzylphthalate             |      | 25     | 2.0 |
| 35 | Benzo(a)anthracene               |      | 25     | 2.0 |
| 36 | bis(2-Ethylhexyl)phthalate       |      | 25     | 2.0 |
| 37 | Chrysene                         |      | 25     | 2.0 |
| 38 | Di-n-octylphthalate              |      | 25     | 2.0 |
| 39 | Benzo(b)fluoranthene             |      | 24     | 2.0 |
| 40 | Benzo(k)fluoranthene             |      | 26     | 2.0 |
| 41 | Benzo(a)pyrene                   |      | 25     | 2.0 |
| 42 | Indeno(1,2,3-cd)pyrene           |      | 24     | 2.0 |
| 43 | Dibenz(a,h)anthracene            |      | 24     | 2.0 |
| 44 | Benzo(g,h,i)perylene             |      | 24     | 2.0 |

## Quantitation Report

(QT Reviewed)

Data File : C:\HPCHEM\1\DATA\NOV2901\CALI25.D  
 Acq On : 29 Nov 2001 4:31 pm  
 Sample : calib 25  
 Misc :  
 MS Integration Params: RTEINT.P  
 Quant Time: Nov 30 14:50 2001

Vial: 3  
 Operator: 209 GALOS  
 Inst : GC/MS 209  
 Multiplr: 1.00

Quant Results File: GCMS1126.RES

Quant Method : C:\HPCHEM\1\METHODS\GCMS1126.M (RTE Integrator)  
 Title : 209 625/TCLP calibration table  
 Last Update : Wed Nov 28 15:06:24 2001  
 Response via : Initial Calibration  
 DataAcq Meth : GCMS1126

| Internal Standards        | R.T.  | QIon | Response | Conc  | Units | Dev(Min) |
|---------------------------|-------|------|----------|-------|-------|----------|
| 1) 1,4-Dichlorobenzene-d4 | 11.70 | 152  | 767415   | 20.00 | ug/l  | 0.00     |
| 15) Naphthalene-d8        | 15.30 | 136  | 2988169  | 20.00 | ug/l  | 0.00     |
| 26) Acenaphthene-d10      | 20.57 | 164  | 1387791  | 20.00 | ug/l  | 0.00     |
| 42) Phenanthrene-d10      | 25.00 | 188  | 1952442  | 20.00 | ug/l  | 0.00     |
| 53) Chrysene-d12          | 33.03 | 240  | 1021146  | 20.00 | ug/l  | 0.00     |
| 59) Perylene-d12          | 37.12 | 264  | 641195   | 20.00 | ug/l  | 0.00     |

## System Monitoring Compounds

|               |       |     |          |      |        |       |
|---------------|-------|-----|----------|------|--------|-------|
| 52) 2,4-DDD   | 29.91 | 235 | 11832    | 0.51 | ug/mL  | -0.16 |
| Spiked Amount | 5.000 |     | Recovery | =    | 10.20% |       |

## Target Compounds

|                                |       |     |          |       |      | Qvalue |
|--------------------------------|-------|-----|----------|-------|------|--------|
| 2) Pyridine                    | 5.36  | 79  | 1348931m | 24.69 | ug/l |        |
| 3) N-nitroso-dimethyl amine    | 5.42  | 74  | 840084   | 24.89 | ug/l | 98     |
| 4) Phenol                      | 10.97 | 94  | 1685029  | 24.90 | ug/l | 98     |
| 5) bis(2-Chloroethyl)ether     | 11.14 | 93  | 1296360  | 24.47 | ug/l | 100    |
| 6) 2-Chlorophenol              | 11.24 | 128 | 1272130  | 24.86 | ug/l | 98     |
| 7) 1,3-Dichlorobenzene         | 11.60 | 146 | 1391332  | 24.93 | ug/l | 100    |
| 8) 1,4-Dichlorobenzene         | 11.74 | 146 | 1405625  | 24.91 | ug/l | 99     |
| 9) 1,2-Dichlorobenzene         | 12.26 | 146 | 1292380  | 24.83 | ug/l | 98     |
| 10) 2-Methylphenol             | 12.55 | 108 | 1182529  | 25.54 | ug/l | 99     |
| 11) 2,2'-oxybis(1-Chloropropan | 12.60 | 45  | 1732774  | 25.69 | ug/l | 98     |
| 12) 3&4-Methylphenol           | 12.96 | 108 | 1204983  | 25.40 | ug/l | 100    |
| 13) N-Nitroso-di-n-propylamine | 12.99 | 70  | 758060   | 25.12 | ug/l | 98     |
| 14) Hexachloroethane           | 13.08 | 117 | 517929   | 25.23 | ug/l | 96     |
| 16) Nitrobenzene               | 13.38 | 77  | 1158626  | 24.88 | ug/l | 99     |
| 17) Isophorone                 | 14.03 | 82  | 2207828  | 24.60 | ug/l | 99     |
| 18) 2-Nitrophenol              | 14.29 | 139 | 686051   | 23.48 | ug/l | 97     |
| 19) 2,4-Dimethylphenol         | 14.45 | 107 | 1096006  | 25.00 | ug/l | 99     |
| 20) bis(2-Chloroethoxy)methane | 14.72 | 93  | 1495393  | 24.81 | ug/l | 99     |
| 21) 2,4-Dichlorophenol         | 14.97 | 162 | 961854   | 24.88 | ug/l | 99     |
| 22) 1,2,4-Trichlorobenzene     | 15.19 | 180 | 1067356  | 24.60 | ug/l | 99     |
| 23) Naphthalene                | 15.36 | 128 | 3596896  | 25.19 | ug/l | 100    |
| 24) Hexachlorobutadiene        | 15.91 | 225 | 527327   | 24.94 | ug/l | 100    |
| 25) 4-Chloro-3-methylphenol    | 17.10 | 107 | 960204   | 24.87 | ug/l | 97     |
| 27) Hexachlorocyclopentadiene  | 18.09 | 237 | 165096   | 16.51 | ug/l | 99     |
| 28) 2,4,6-Trichlorophenol      | 18.36 | 196 | 593915   | 24.84 | ug/l | 99     |
| 29) 2,4,5-Trichlorophenol      | 18.47 | 196 | 641200   | 24.44 | ug/l | 100    |
| 30) 2-Chloronaphthalene        | 18.85 | 162 | 2015683  | 25.46 | ug/l | 99     |
| 31) Dimethylphthalate          | 19.96 | 163 | 2302504  | 24.97 | ug/l | 100    |
| 32) 2,6-Dinitrotoluene         | 20.17 | 165 | 534650m  | 24.53 | ug/l |        |

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

CALI25.D GCMS1126.M

Fri Nov 30 14:51:03 2001

Page 1

## Quantitation Report (QT Reviewed)

Data File : C:\HPCHEM\1\DATA\NOV2901\CALI25.D  
 Acq On : 29 Nov 2001 4:31 pm  
 Sample : calib 25  
 Misc :  
 MS Integration Params: RTEINT.P  
 Quant Time: Nov 30 14:50 2001

Vial: 3  
 Operator: 209 GALOS  
 Inst : GC/MS 209  
 Multiplr: 1.00

Quant Results File: GCMS1126.RES

Quant Method : C:\HPCHEM\1\METHODS\GCMS1126.M (RTE Integrator)  
 Title : 209 625/TCLP calibration table  
 Last Update : Wed Nov 28 15:06:24 2001  
 Response via : Initial Calibration  
 DataAcq Meth : GCMS1126

| Compound                        | R.T.  | QIon | Response | Conc Unit    | Qvalue |
|---------------------------------|-------|------|----------|--------------|--------|
| 33) Acenaphthylene              | 20.11 | 152  | 2818424  | 25.33 ug/l   | 99     |
| 34) Acenaphthene                | 20.68 | 153  | 2013576  | 25.31 ug/l   | 100    |
| 35) 2,4-Dinitrophenol           | 20.88 | 184  | 107105   | 12.36 ug/l   | 91     |
| 36) 4-Nitrophenol               | 21.15 | 65   | 204793   | 22.60 ug/l   | 94     |
| 37) 2,4-Dinitrotoluene          | 21.34 | 165  | 668576   | 24.13 ug/l   | 96     |
| 38) Diethylphthalate            | 22.10 | 149  | 2147351  | 25.16 ug/l   | 99     |
| 39) 4-Chlorophenyl-phenylether  | 22.22 | 204  | 949379   | 25.48 ug/l   | 99     |
| 40) Fluorene                    | 22.19 | 166  | 2055384  | 25.18 ug/l   | 100    |
| 41) Azobenzene/ 1,2-Diphenylhyd | 22.69 | 77   | 2204595  | 25.00 ug/L   | 98     |
| 43) 4,6-Dinitro-2-methylphenol  | 22.55 | 198  | 256791   | 18.80 ug/l   | 99     |
| 44) N-Nitrosodiphenylamine      | 22.62 | 168  | 916238   | 26.43 ug/l   | 100    |
| 45) 4-Bromophenyl-phenylether   | 23.68 | 248  | 484770   | 25.24 ug/l   | 98     |
| 46) Hexachlorobenzene           | 24.10 | 284  | 520403   | 24.64 ug/l   | 98     |
| 47) Pentachlorophenol           | 24.67 | 266  | 216935   | 21.71 ug/l   | 99     |
| 48) Phenanthrene                | 25.06 | 178  | 2596002  | 25.11 ug/l   | 99     |
| 49) Anthracene                  | 25.20 | 178  | 2576560  | 25.00 ug/l   | 99     |
| 50) Di-n-butylphthalate         | 27.01 | 149  | 3279920  | 23.84 ug/l   | 100    |
| 51) Fluoranthene                | 28.67 | 202  | 2351786  | 23.56 ug/l   | 96     |
| 54) Pyrene                      | 29.34 | 202  | 2367947  | 26.87 ug/l   | 99     |
| 55) Butylbenzylphthalate        | 31.51 | 149  | 1061711  | 23.06 ug/l   | 93     |
| 56) Benzo(a)anthracene          | 32.98 | 228  | 1496286  | 24.59 ug/l   | 99     |
| 57) Bis(2-ethylhexyl)phthalate  | 33.31 | 149  | 1274810  | 20.49 ug/l   | 100    |
| 58) Chrysene                    | 33.11 | 228  | 1449984  | 24.30 ug/l   | 99     |
| 60) Di-n-Octylphthalate         | 35.04 | 149  | 1788301  | 20.99 ug/l # | 97     |
| 61) Benzo(b)fluoranthene        | 36.05 | 252  | 1199228  | 24.86 ug/l   | 96     |
| 62) Benzo(k)fluoranthene        | 36.12 | 252  | 1276442  | 25.78 ug/l   | 99     |
| 63) Benzo(a)pyrene              | 36.94 | 252  | 1025558  | 24.45 ug/l   | 99     |
| 64) Indeno(1,2,3-cd)pyrene      | 40.80 | 276  | 1014319  | 24.19 ug/l   | 99     |
| 65) Dibenz(a,h)anthracene       | 40.86 | 278  | 811035   | 24.70 ug/l   | 99     |
| 66) Benzo(g,h,i)perylene        | 41.90 | 276  | 861653   | 24.05 ug/l   | 97     |

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

CALI25.D GCMS1126.M Fri Nov 30 14:51:07 2001

Page 2

## Quantitation Report

Data File : C:\HPCHEM\1\DATA\NOV2901\CALI25.D  
Acq On : 29 Nov 2001 4:31 pm  
Sample : calib 25  
Misc :  
MS Integration Params: RTEINT.P  
Quant Time: Nov 30 14:50 2001 Qu.

Vial: 3  
Operator: 209 GALOS  
Inst : GC/MS 209  
Multiplr: 1.00

Quant Results File: GCMS1126.RES

```
Method      : C:\HPCHEM\1\METHODS\GCMS1126.M (RTE Integrator)
Title       : 209 625/TCLP calibration table
Last Update : Wed Nov 28 15:06:24 2001
Response via : Initial Calibration
```

