User: Galos, Marie
Westchester Dept. of Labs & Research
Date: 10/16/2001 Time: 09:14:23

Sample: AD23326
Analysis: $GCMS GCMS Scan For Unknowns
Units: ug
Started: 10/16/01 09:13 Ended: 10/16/01 09:13 Analyst: MG
Page: 1

| | Component Name | Viol | Result |
|---|---|---|---|
| 1 | Other unknown compounds | | see comment |

**Comments for Sample AD23326 - Analysis $GCMS**

Westchester Dept. of Labs & Research
User: Galos, Marie
Date: 10-16-2001 Time: 09:14:37

A scan of the extract prepared for the 508 analysis, from 35-550 amu using a mass spectrometer, does not reveal the presence of non-target compounds.

Quantitation Report (QT Reviewed)

Data File : C:\HPCHEM\1\DATA\OCT0901\23326.D Vial: 13
Acq On : 10 Oct 2001 1:04 am Operator: 209 GALOS
Sample : gc/ms unknown Inst : GC/MS 209
Misc : Multiplr: 1.00
MS Integration Params: RTEINT.P
Quant Time: Oct 11 10:28 2001 Quant Results File: NP091101.RES

Quant Method : C:\HPCHEM\1\METHODS\NP091101.M (RTE Integrator)
Title : 209 625/TCLP calibration table 7/16/01
Last Update : Thu Sep 13 12:47:04 2001
Response via : Initial Calibration
DataAcq Meth : NP091101

| Internal Standards | R.T. | QIon | Response | Conc | Units | Dev(Min) |
|---|---|---|---|---|---|---|
| 1) 1,4-Dichlorobenzene-d4 | 11.76 | 152 | 350806 | 20.00 | ug/l | -0.14 |
| 19) Naphthalene-d8 | 15.37 | 136 | 1398665 | 20.00 | ug/l | -0.15 |
| 31) Acenaphthene-d10 | 20.66 | 164 | 641090 | 20.00 | ug/l | -0.15 |
| 49) Phenanthrene-d10 | 25.09 | 188 | 899802 | 20.00 | ug/l | -0.15 |
| 59) Chrysene-d12 | 33.12 | 240 | 592191 | 20.00 | ug/l | -0.16 |
| 68) Perylene-d12 | 37.23 | 264 | 440679 | 20.00 | ug/l | -0.19 |
| **System Monitoring Compounds** | | | | | | |
| 4) 2-Fluorophenol | 0.00 | 112 | 0 | 0.00 | ug/l | |
| Spiked Amount 75.000 | | | Recovery = | 0.00% | | |
| 5) Phenol-d5 | 0.00 | 99 | 0 | 0.00 | ug/l | |
| Spiked Amount 75.000 | | | Recovery = | 0.00% | | |
| 6) 2-Chlorophenol-d4 | 0.00 | 132 | 0 | 0.00 | ug/l | |
| Spiked Amount 75.000 | | | Recovery = | 0.00% | | |
| 7) 1,2-Dichlorobenzene-d4 | 12.29 | 152 | 3569 | 0.21 | ug/l | -0.14 |
| Spiked Amount 50.000 | | | Recovery = | 0.42% | | |
| 20) Nitrobenzene-d5 | 0.00 | 82 | 0 | 0.00 | ug/l | |
| Spiked Amount 50.000 | | | Recovery = | 0.00% | | |
| 32) 2-Fluorobiphenyl | 18.81 | 172 | 1280 | 0.03 | ug/l | 0.00 |
| Spiked Amount 50.000 | | | Recovery = | 0.06% | | |
| 33) 2,4,6-Tribromophenol | 0.00 | 330 | 0 | 0.00 | ug/l | |
| Spiked Amount 75.000 | | | Recovery = | 0.00% | | |
| 62) Terphenyl-d14 | 0.00 | 244 | 0 | 0.00 | ug/l | |
| Spiked Amount 50.000 | | | Recovery = | 0.00% | | |
| **Target Compounds** | | | | | | Qvalue |
| 2) Pyridine | 0.00 | 79 | 0 | N.D. | | |
| 3) N-nitroso-dimethyl amine | 0.00 | 74 | 0 | N.D. | | |
| 8) Phenol | 0.00 | 94 | 0 | N.D. | | |
| 9) bis(2-Chloroethyl)ether | 0.00 | 93 | 0 | N.D. | | |
| 10) 2-Chlorophenol | 0.00 | 128 | 0 | N.D. | | |
| 11) 1,3-Dichlorobenzene | 0.00 | 146 | 0 | N.D. | | |
| 12) 1,4-Dichlorobenzene | 0.00 | 146 | 0 | N.D. | | |
| 13) 1,2-Dichlorobenzene | 0.00 | 146 | 0 | N.D. | | |
| 14) 2-Methylphenol | 0.00 | 108 | 0 | N.D. | | |
| 15) 2,2'-oxybis(1-Chloropropan | 0.00 | 45 | 0 | N.D. | | |
| 16) 3&4-Methylphenol | 0.00 | 108 | 0 | N.D. | | |
| 17) N-Nitroso-di-n-propylamine | 0.00 | 70 | 0 | N.D. | | |
| 18) Hexachloroethane | 0.00 | 117 | 0 | N.D. | | |
| 21) Nitrobenzene | 0.00 | 77 | 0 | N.D. | | |
| 22) Isophorone | 0.00 | 82 | 0 | N.D. | | |

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

Quantitation Report (QT Reviewed)

Data File : C:\HPCHEM\1\DATA\OCT0901\23326.D Vial: 13
Acq On : 10 Oct 2001 1:04 am Operator: 209 GALOS
Sample : gc/ms unknown Inst : GC/MS 209
Misc : Multiplr: 1.00
MS Integration Params: RTEINT.P
Quant Time: Oct 11 10:28 2001 Quant Results File: NP091101.RES

Quant Method : C:\HPCHEM\1\METHODS\NP091101.M (RTE Integrator)
Title : 209 625/TCLP calibration table 7/16/01
Last Update : Thu Sep 13 12:47:04 2001
Response via : Initial Calibration
DataAcq Meth : NP091101

| | Compound | R.T. | QIon | Response | Conc Unit | Qvalue |
|---|---|---|---|---|---|---|
| 23) | 2-Nitrophenol | 0.00 | 139 | 0 | N.D. | |
| 24) | 2,4-Dimethylphenol | 0.00 | 107 | 0 | N.D. | |
| 25) | bis(2-Chloroethoxy)methane | 0.00 | 93 | 0 | N.D. | |
| 26) | 2,4-Dichlorophenol | 0.00 | 162 | 0 | N.D. | |
| 27) | 1,2,4-Trichlorobenzene | 0.00 | 180 | 0 | N.D. | |
| 28) | Naphthalene | 0.00 | 128 | 0 | N.D. | |
| 29) | Hexachlorobutadiene | 0.00 | 225 | 0 | N.D. | |
| 30) | 4-Chloro-3-methylphenol | 0.00 | 107 | 0 | N.D. | |
| 34) | Hexachlorocyclopentadiene | 0.00 | 237 | 0 | N.D. | |
| 35) | 2,4,6-Trichlorophenol | 0.00 | 196 | 0 | N.D. | |
| 36) | 2,4,5-Trichlorophenol | 0.00 | 196 | 0 | N.D. | |
| 37) | 2-Chloronaphthalene | 0.00 | 162 | 0 | N.D. | |
| 38) | Dimethylphthalate | 0.00 | 163 | 0 | N.D. | |
| 39) | 2,6-Dinitrotoluene | 0.00 | 165 | 0 | N.D. | d |
| 40) | Acenaphthylene | 0.00 | 152 | 0 | N.D. | |
| 41) | Acenaphthene | 0.00 | 153 | 0 | N.D. | |
| 42) | 2,4-Dinitrophenol | 0.00 | 184 | 0 | N.D. | |
| 43) | 4-Nitrophenol | 0.00 | 65 | 0 | N.D. | |
| 44) | 2,4-Dinitrotoluene | 0.00 | 165 | 0 | N.D. | |
| 45) | Diethylphthalate | 0.00 | 149 | 0 | N.D. | |
| 46) | 4-Chlorophenyl-phenylether | 0.00 | 204 | 0 | N.D. | |
| 47) | Fluorene | 0.00 | 166 | 0 | N.D. | |
| 48) | Azobenzen/ 1,2-Diphenylhyd | 0.00 | 77 | 0 | N.D. | |
| 50) | 4,6-Dinitro-2-methylphenol | 0.00 | 198 | 0 | N.D. | |
| 51) | N-Nitrosodiphenylamine | 0.00 | 168 | 0 | N.D. | |
| 52) | 4-Bromophenyl-phenylether | 0.00 | 248 | 0 | N.D. | |
| 53) | Hexachlorobenzene | 0.00 | 284 | 0 | N.D. | |
| 54) | Pentachlorophenol | 0.00 | 266 | 0 | N.D. | |
| 55) | Phenanthrene | 25.08 | 178 | 101 | N.D. | |
| 56) | Anthracene | 0.00 | 178 | 0 | N.D. | |
| 57) | Di-n-butylphthalate | 27.09 | 149 | 2341 | N.D. | |
| 58) | Fluoranthene | 0.00 | 202 | 0 | N.D. | |
| 60) | 3,3'-Dichlorobenzidine | 0.00 | 252 | 0 | N.D. | |
| 61) | Benzidine | 0.00 | 184 | 0 | N.D. | |
| 63) | Pyrene | 0.00 | 202 | 0 | N.D. | |
| 64) | Butylbenzylphthalate | 0.00 | 149 | 0 | N.D. | |
| 65) | Benzo(a)anthracene | 33.13 | 228 | 1288 | N.D. | |
| 66) | Bis(2-ethylhexyl)phthalate | 33.38 | 149 | 3535 | N.D. | |
| 67) | Chrysene | 33.20 | 228 | 486 | N.D. | |
| 69) | Di-n-Octylphthalate | 0.00 | 149 | 0 | N.D. | |
| 70) | Benzo(b)fluoranthene | 0.00 | 252 | 0 | N.D. | |

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

Quantitation Report (QT Reviewed)

Data File : C:\HPCHEM\1\DATA\OCT0901\23326.D Vial: 13
Acq On : 10 Oct 2001 1:04 am Operator: 209 GALOS
Sample : gc/ms unknown Inst : GC/MS 209
Misc : Multiplr: 1.00
MS Integration Params: RTEINT.P
Quant Time: Oct 11 10:28 2001 Quant Results File: NP091101.RES

Quant Method : C:\HPCHEM\1\METHODS\NP091101.M (RTE Integrator)
Title : 209 625/TCLP calibration table 7/16/01
Last Update : Thu Sep 13 12:47:04 2001
Response via : Initial Calibration
DataAcq Meth : NP091101

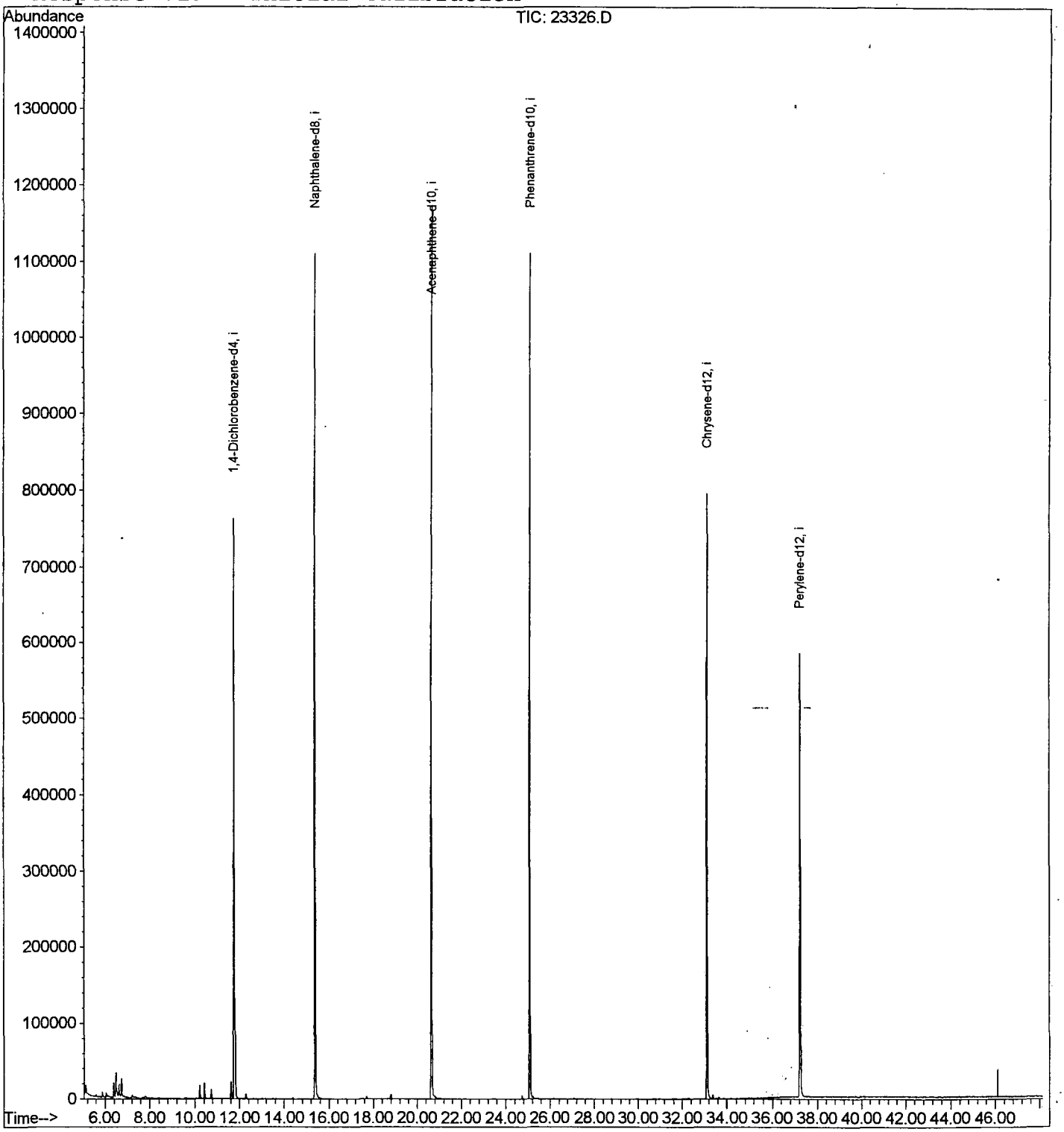
| | Compound | R.T. | QIon | Response | Conc Unit | Qvalue |
|---|---|---|---|---|---|---|
| 71) | Benzo(k)fluoranthene | 0.00 | 252 | 0 | N.D. | |
| 72) | Benzo(a)pyrene | 37.23 | 252 | 1584 | N.D. | |
| 73) | Indeno(1,2,3-cd)pyrene | 0.00 | 276 | 0 | N.D. | |
| 74) | Dibenz(a,h)anthracene | 41.04 | 278 | 540 | N.D. | |
| 75) | Benzo(g,h,i)perylene | 42.11 | 276 | 1018 | N.D. | |

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

# Quantitation Report

Data File : C:\HPCHEM\1\DATA\OCT0901\23326.D Vial: 13
Acq On : 10 Oct 2001 1:04 am Operator: 209 GALOS
Sample : gc/ms unknown Inst : GC/MS 209
Misc : Multiplr: 1.00
MS Integration Params: RTEINT.P
Quant Time: Oct 11 10:28 2001 Quant Results File: NP091101.RES

Method : C:\HPCHEM\1\METHODS\NP091101.M (RTE Integrator)
Title : 209 625/TCLP calibration table 7/16/01
Last Update : Thu Sep 13 12:47:04 2001
Response via : Initial Calibration



# LSC Area Percent Report

Data File : C:\HPCHEM\1\DATA\OCT0901\23326.D  Vial: 13
Acq On : 10 Oct 2001 1:04 am  Operator: 209 GALOS
Sample : gc/ms unknown  Inst : GC/MS 209
Misc :  Multiplr: 1.00
MS Integration Params: LSCINT.P

Method : C:\HPCHEM\1\METHODS\NP091101.M (RTE Integrator)
Title : 209 625/TCLP calibration table 7/16/01
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

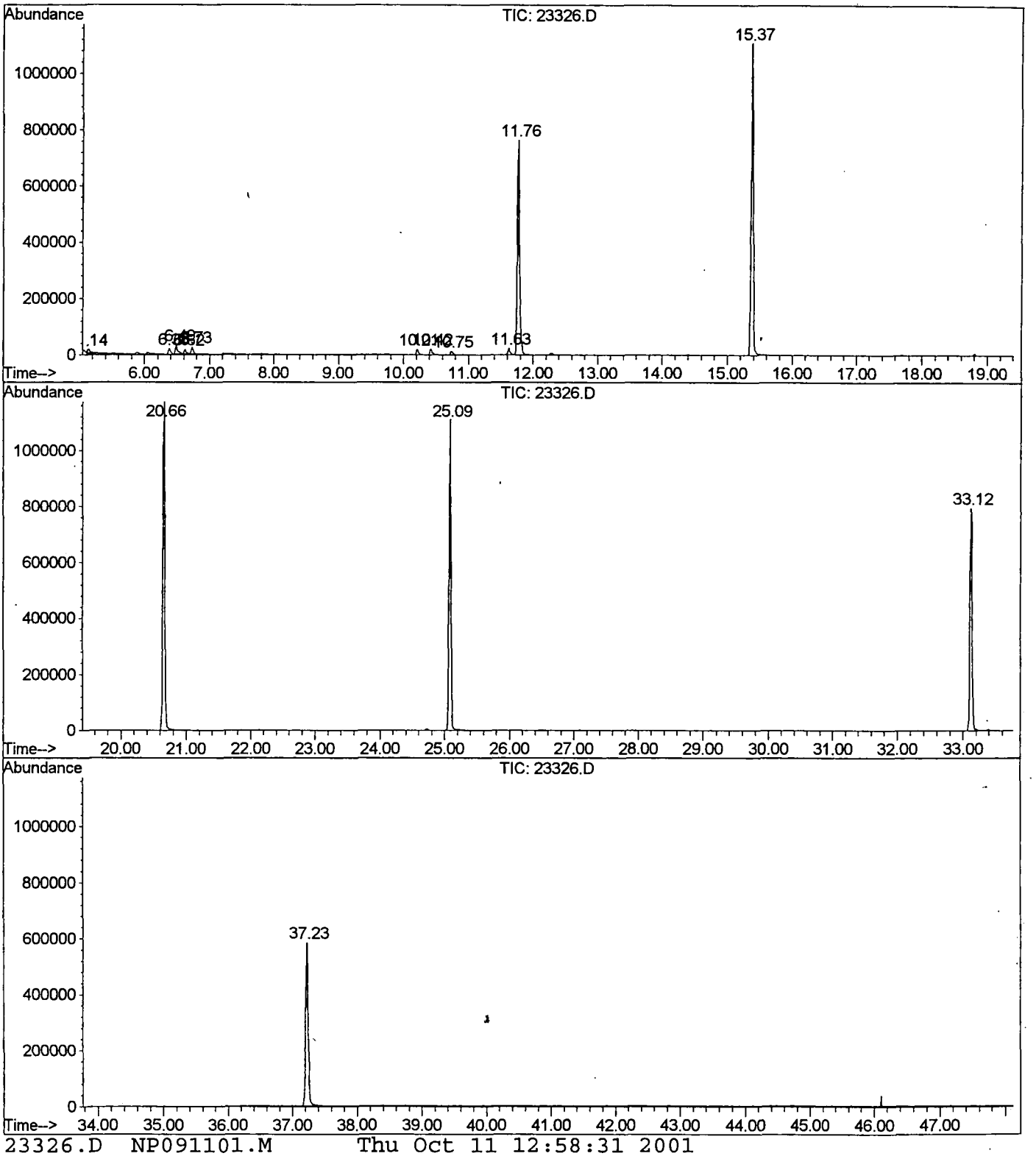
| peak # | R.T. min | first scan | max scan | last scan | PK TY | peak height | peak area | peak % max. | % of total |
|---|---|---|---|---|---|---|---|---|---|
| 1 | 5.136 | 8 | 10 | 23 | rVB | 12083 | 31635 | 1.23% | 0.241% |
| 2 | 6.385 | 141 | 146 | 152 | rBV | 19611 | 41168 | 1.60% | 0.313% |
| 3 | 6.486 | 153 | 157 | 167 | rVV2 | 31793 | 82182 | 3.19% | 0.625% |
| 4 | 6.623 | 168 | 172 | 177 | rVV | 15961 | 35453 | 1.38% | 0.270% |
| 5 | 6.734 | 180 | 184 | 192 | rVB | 23699 | 50824 | 1.97% | 0.387% |
| 6 | 10.213 | 559 | 563 | 577 | rVB | 18046 | 40208 | 1.56% | 0.306% |
| 7 | 10.424 | 582 | 586 | 595 | rBV | 20509 | 46674 | 1.81% | 0.355% |
| 8 | 10.745 | 616 | 621 | 628 | rBV | 12118 | 27276 | 1.06% | 0.208% |
| 9 | 11.627 | 711 | 717 | 727 | rBV | 22643 | 50465 | 1.96% | 0.384% |
| 10 | 11.764 | 727 | 732 | 746 | rBV | 762115 | 1830351 | 71.04% | 13.928% |
| 11 | 15.372 | 1119 | 1125 | 1143 | rBV | 1109426 | 2549671 | 98.96% | 19.402% |
| 12 | 20.660 | 1694 | 1701 | 1719 | rBV | 1173430 | 2576586 | 100.00% | 19.607% |
| 13 | 25.085 | 2176 | 2183 | 2194 | rBV | 1110478 | 2322543 | 90.14% | 17.674% |
| 14 | 33.118 | 3051 | 3058 | 3074 | rBV2 | 794825 | 1841725 | 71.48% | 14.015% |
| 15 | 37.231 | 3498 | 3506 | 3520 | rBV2 | 582581 | 1614301 | 62.65% | 12.284% |

Sum of corrected areas: 13141062

23326.D NP091101.M Thu Oct 11 12:58:29 2001

# LSC Report - Integrated Chromatogram

File : C:\HPCHEM\1\DATA\OCT0901\23326.D
Operator : 209 GALOS
Acquired : 10 Oct 2001 1:04 am using AcqMethod NP091101
Instrument : GC/MS 209
Sample Name: gc/ms unknown
Misc Info :
Vial Number: 13
Quant File :NP091101.RES (RTE Integrator)




23326.D NP091101.M Thu Oct 11 12:58:31 2001

# Library Search Compound Report

Data File : C:\HPCHEM\1\DATA\OCT0901\23326.D
Acq On : 10 Oct 2001 1:04 am
Sample : gc/ms unknown
Misc :
MS Integration Params: LSCINT.P

Vial: 13
Operator: 209 GALOS
Inst : GC/MS 209
Multiplr: 1.00

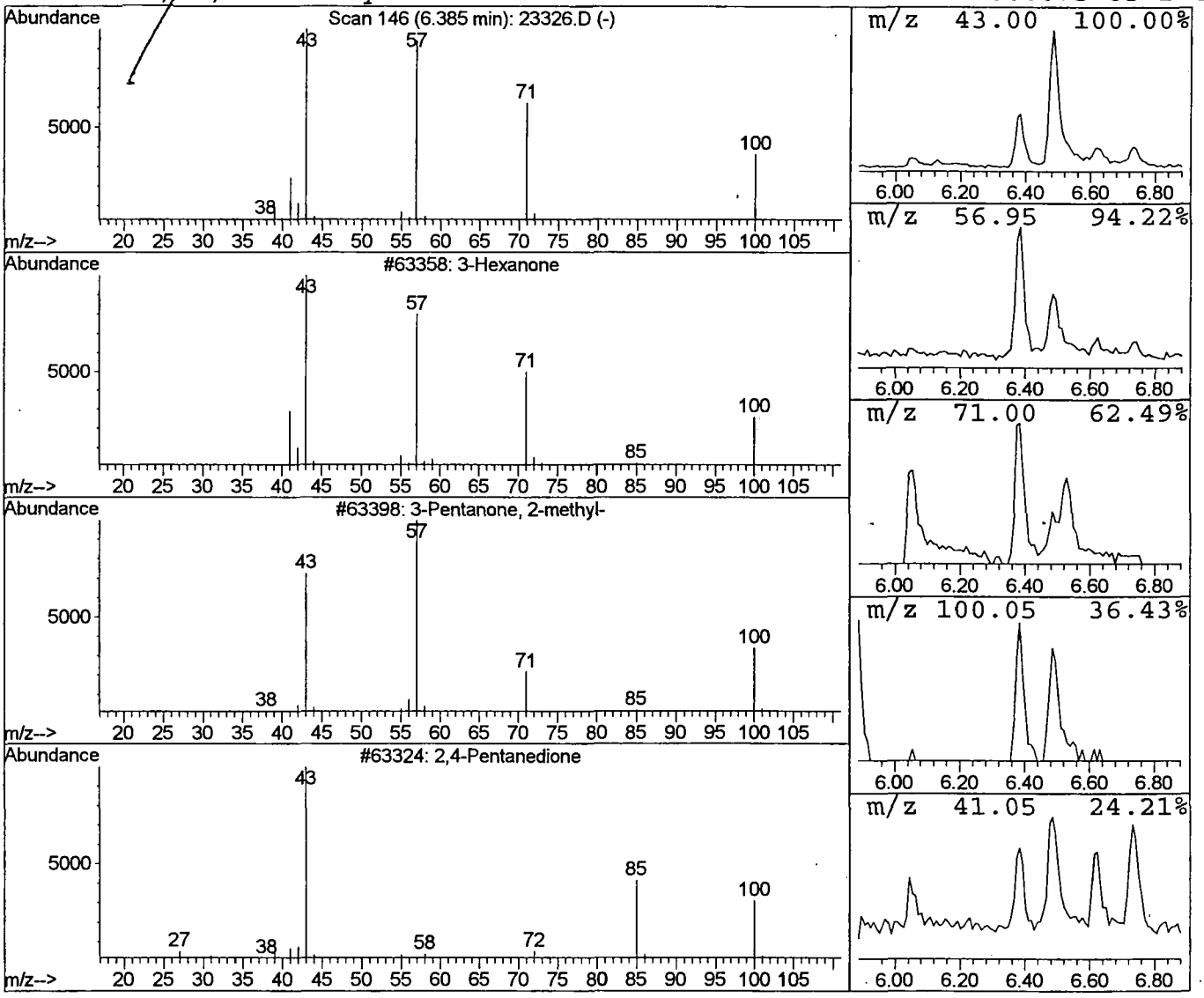
Quant Method : C:\HPCHEM\1\METHODS\NP091101.M (RTE Integrator)
Title : 209 625/TCLP calibration table 7/16/01
Library : C:\DATABASE\NBS75K.L

**********

Peak Number 1 3-Hexanone Concentration Rank 5

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 6.38 | 0.45 ug/l | 41168 | 1,4-Dichlorobenzene-d4 | 11.76 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | 3-Hexanone | 100 | C6H12O | 000589-38-8 | 91 |
| 2 | 3-Pentanone, 2-methyl- | 100 | C6H12O | 000565-69-5 | 53 |
| 3 | 2,4-Pentanedione | 100 | C5H8O2 | 000123-54-6 | 43 |
| 4 | Butane, 2,2-dimethyl- | 86 | C6H14 | 000075-83-2 | 42 |



# Library Search Compound Report

Data File : C:\HPCHEM\1\DATA\OCT0901\23326.D
Acq On : 10 Oct 2001 1:04 am
Sample : gc/ms unknown
Misc :
MS Integration Params: LSCINT.P

Vial: 13
Operator: 209 GALOS
Inst : GC/MS 209
Multiplr: 1.00

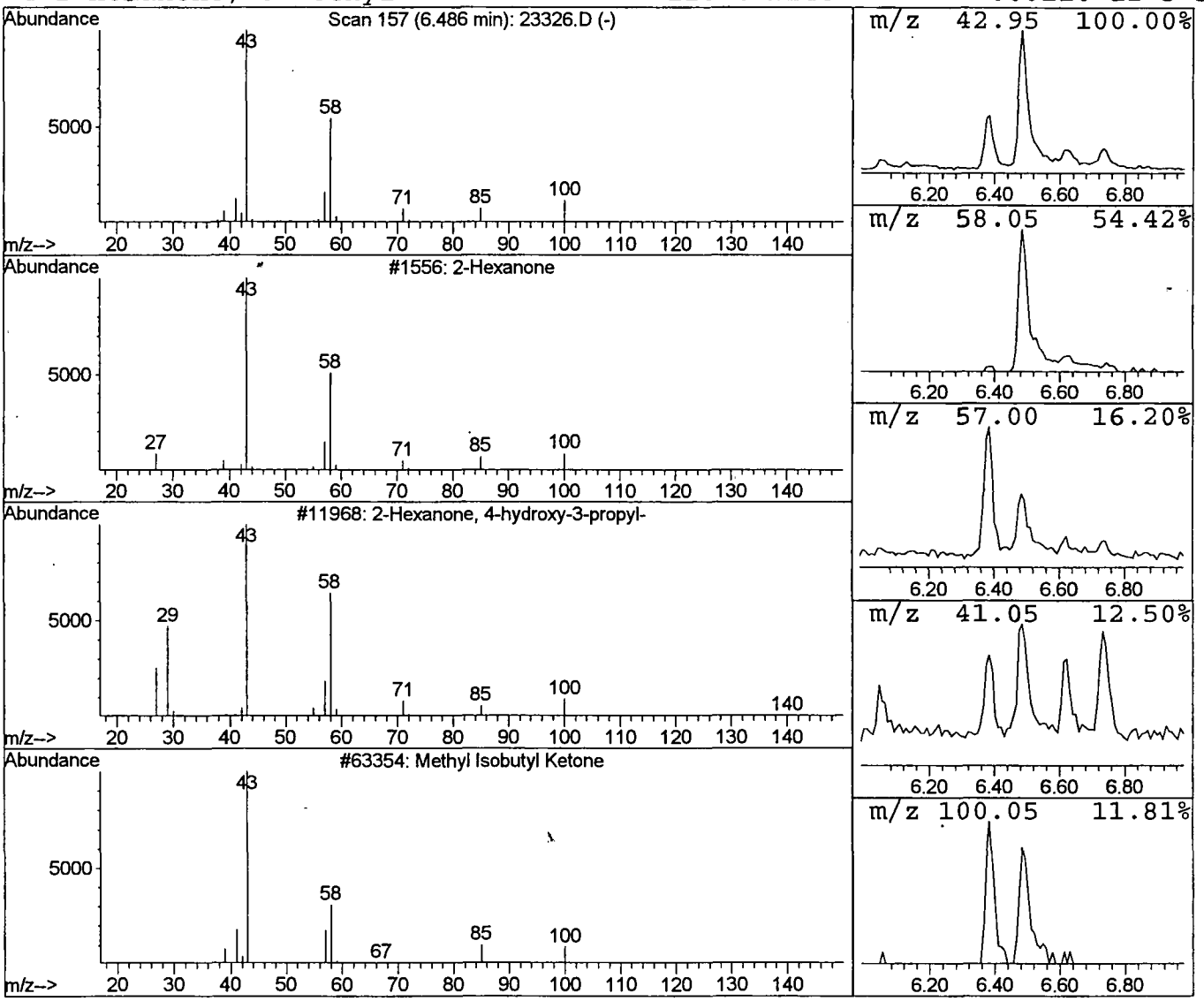
Quant Method : C:\HPCHEM\1\METHODS\NP091101.M (RTE Integrator)
Title : 209 625/TCLP calibration table 7/16/01
Library : C:\DATABASE\NBS75K.L

**********

Peak Number 2 2-Hexanone Concentration Rank 1

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 6.49 | 0.90 ug/l | 82182 | 1,4-Dichlorobenzene-d4 | 11.76 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | 2-Hexanone | 100 | C6H12O | 000591-78-6 | 91 |
| 2 | 2-Hexanone, 4-hydroxy-3-propyl- | 158 | C9H18O2 | 062338-17-4 | 83 |
| 3 | Methyl Isobutyl Ketone | 100 | C6H12O | 000108-10-1 | 53 |
| 4 | 2-Hexanone, 5-methyl- | 114 | C7H14O | 000110-12-3 | 50 |



# Library Search Compound Report

Data File : C:\HPCHEM\1\DATA\OCT0901\23326.D
Acq On : 10 Oct 2001 1:04 am
Sample : gc/ms unknown
Misc :
MS Integration Params: LSCINT.P

Vial: 13
Operator: 209 GALOS
Inst : GC/MS 209
Multiplr: 1.00

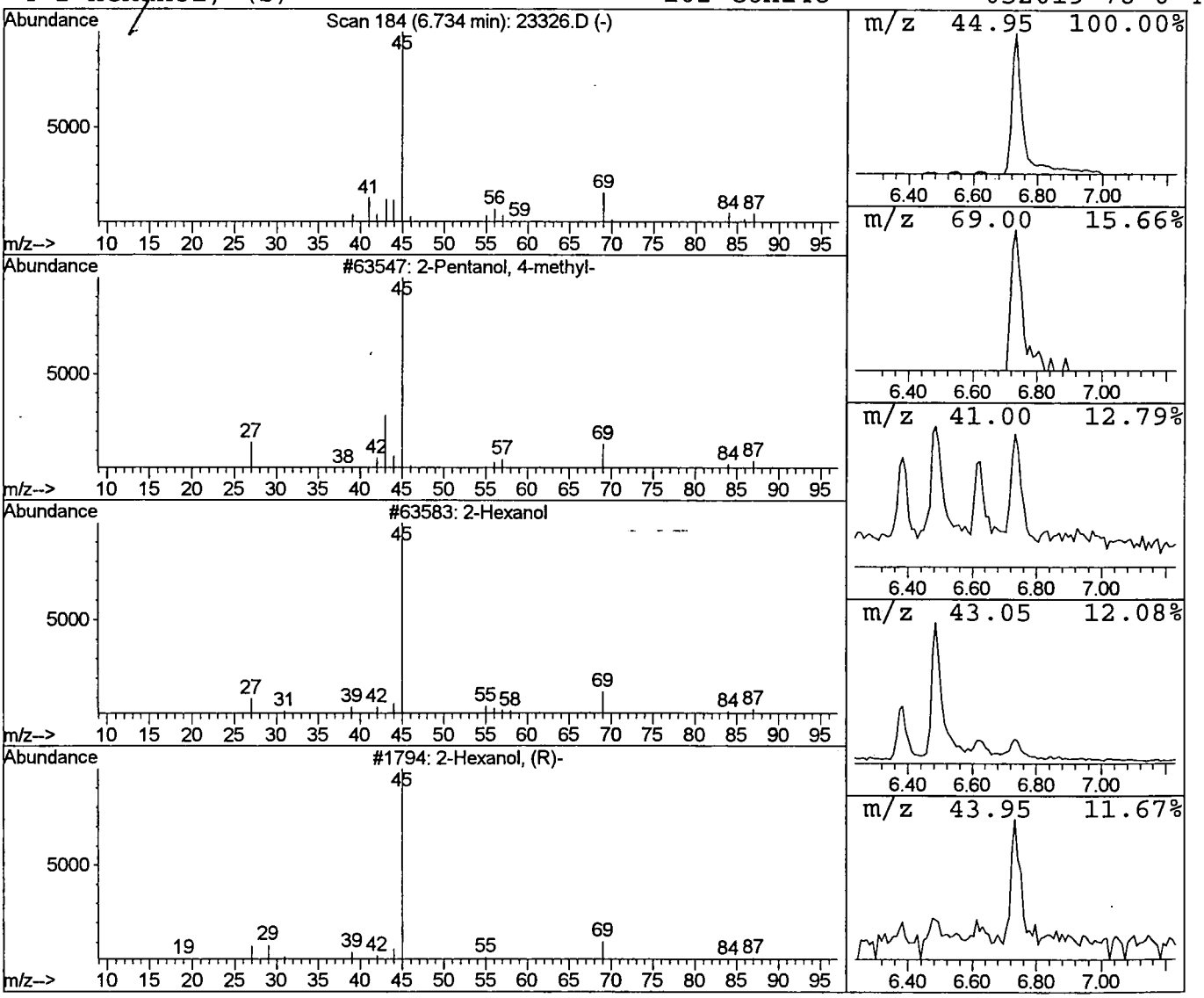
Quant Method : C:\HPCHEM\1\METHODS\NP091101.M (RTE Integrator)
Title : 209 625/TCLP calibration table 7/16/01
Library : C:\DATABASE\NBS75K.L

**********

Peak Number 3 2-Pentanol, 4-methyl- Concentration Rank 2

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 6.73 | 0.56 ug/l | 50824 | 1,4-Dichlorobenzene-d4 | 11.76 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | 2-Pentanol, 4-methyl- | 102 | C6H14O | 000108-11-2 | 83 |
| 2 | 2-Hexanol | 102 | C6H14O | 000626-93-7 | 74 |
| 3 | 2-Hexanol, (R)- | 102 | C6H14O | 026549-24-6 | 74 |
| 4 | 2-Hexanol, (S)- | 102 | C6H14O | 052019-78-0 | 43 |



# Library Search Compound Report

Data File : C:\HPCHEM\1\DATA\OCT0901\23326.D
Acq On : 10 Oct 2001 1:04 am
Sample : gc/ms unknown
Misc :
MS Integration Params: LSCINT.P

Vial: 13
Operator: 209 GALOS
Inst : GC/MS 209
Multiplr: 1.00

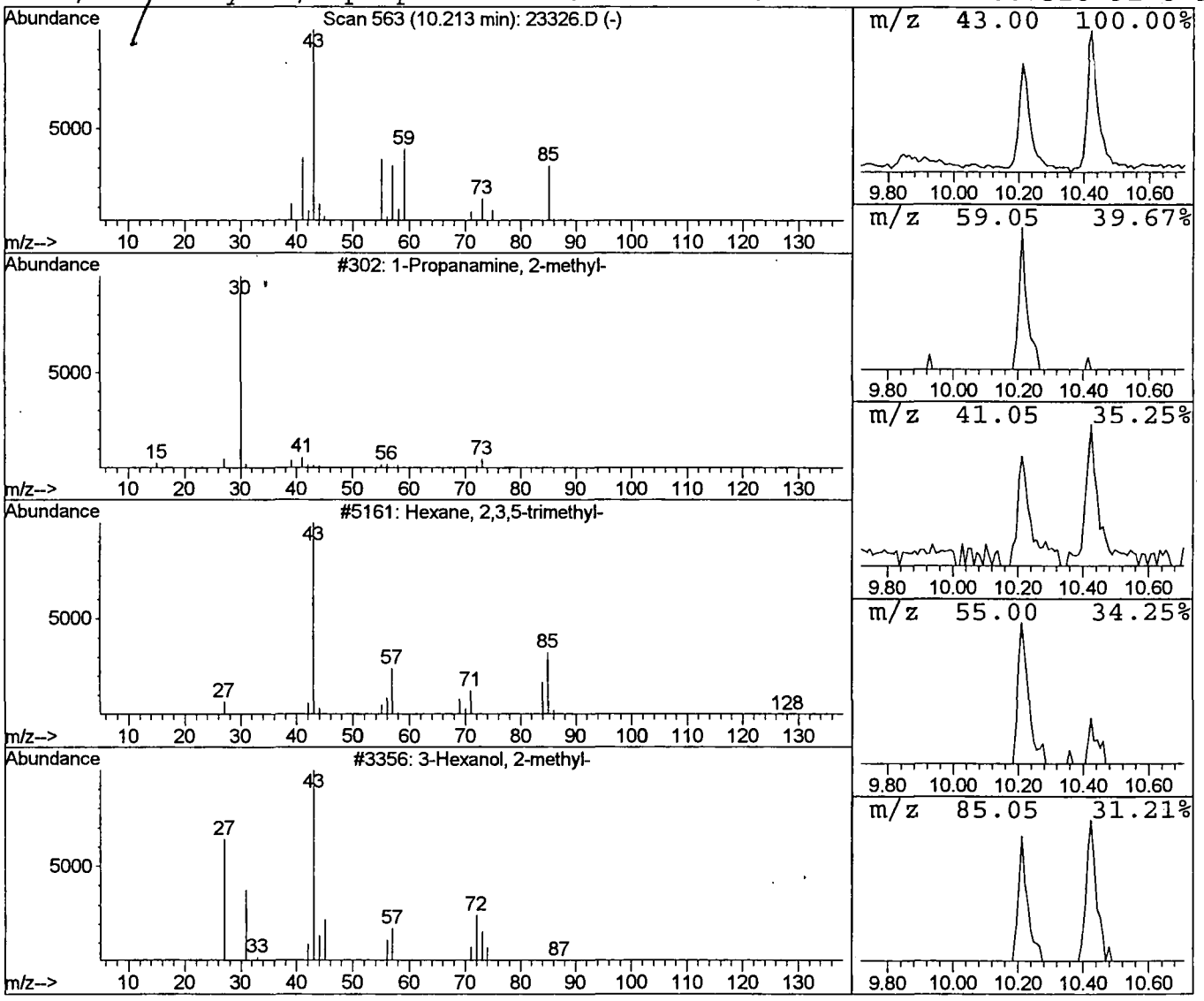
Quant Method : C:\HPCHEM\1\METHODS\NP091101.M (RTE Integrator)
Title : 209 625/TCLP calibration table 7/16/01
Library : C:\DATABASE\NBS75K.L

**********

Peak Number 4 1-Propanamine, 2-methyl- Concentration Rank 6

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 10.21 | 0.44 ug/l | 40208 | 1,4-Dichlorobenzene-d4 | 11.76 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | 1-Propanamine, 2-methyl- | 73 | C4H11N | 000078-81-9 | 9 |
| 2 | Hexane, 2,3,5-trimethyl- | 128 | C9H20 | 001069-53-0 | 9 |
| 3 | 3-Hexanol, 2-methyl- | 116 | C7H16O | 000617-29-8 | 9 |
| 4 | 2,2-Dimethyl-1,3-propanediamine | 102 | C5H14N2 | 007328-91-8 | 9 |



# Library Search Compound Report

Data File : C:\HPCHEM\1\DATA\OCT0901\23326.D
Acq On : 10 Oct 2001 1:04 am
Sample : gc/ms unknown
Misc :
MS Integration Params: LSCINT.P

Vial: 13
Operator: 209 GALOS
Inst : GC/MS 209
Multiplr: 1.00

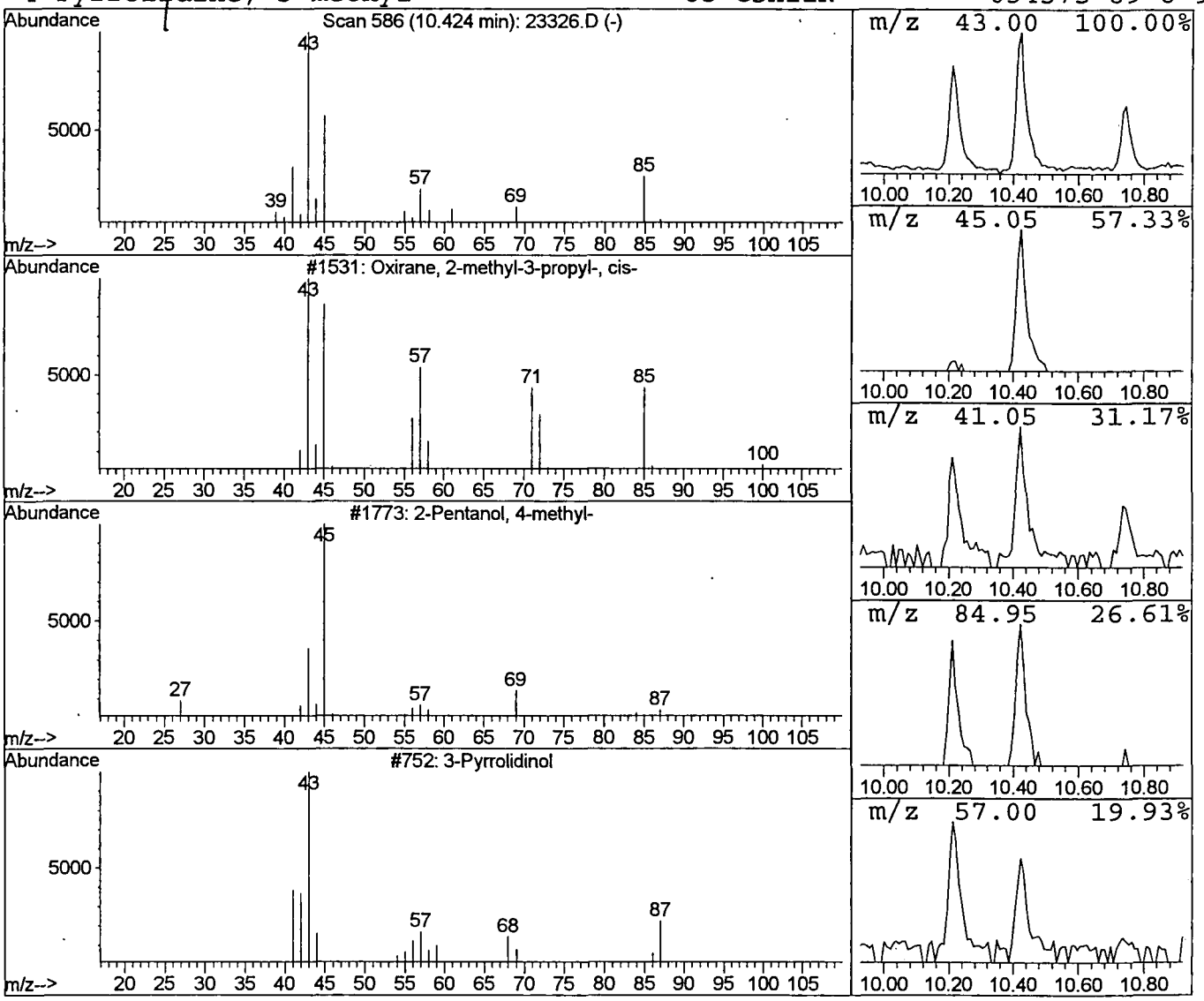
Quant Method : C:\HPCHEM\1\METHODS\NP091101.M (RTE Integrator)
Title : 209 625/TCLP calibration table 7/16/01
Library : C:\DATABASE\NBS75K.L

**********

Peak Number 5 Oxirane, 2-methyl-3-propyl-, c Concentration Rank 4

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 10.42 | 0.51 ug/l | 46674 | 1,4-Dichlorobenzene-d4 | 11.76 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Oxirane, 2-methyl-3-propyl-, cis- | 100 | C6H12O | 006124-90-9 | 33 |
| 2 | 2-Pentanol, 4-methyl- | 102 | C6H14O | 000108-11-2 | 25 |
| 3 | 3-Pyrrolidinol | 87 | C4H9NO | 040499-83-0 | 12 |
| 4 | Pyrrolidine, 3-methyl- | 85 | C5H11N | 034375-89-8 | 9 |



# Library Search Compound Report

Data File : C:\HPCHEM\1\DATA\OCT0901\23326.D
Acq On : 10 Oct 2001 1:04 am
Sample : gc/ms unknown
Misc :
MS Integration Params: LSCINT.P

Vial: 13
Operator: 209 GALOS
Inst : GC/MS 209
Multiplr: 1.00

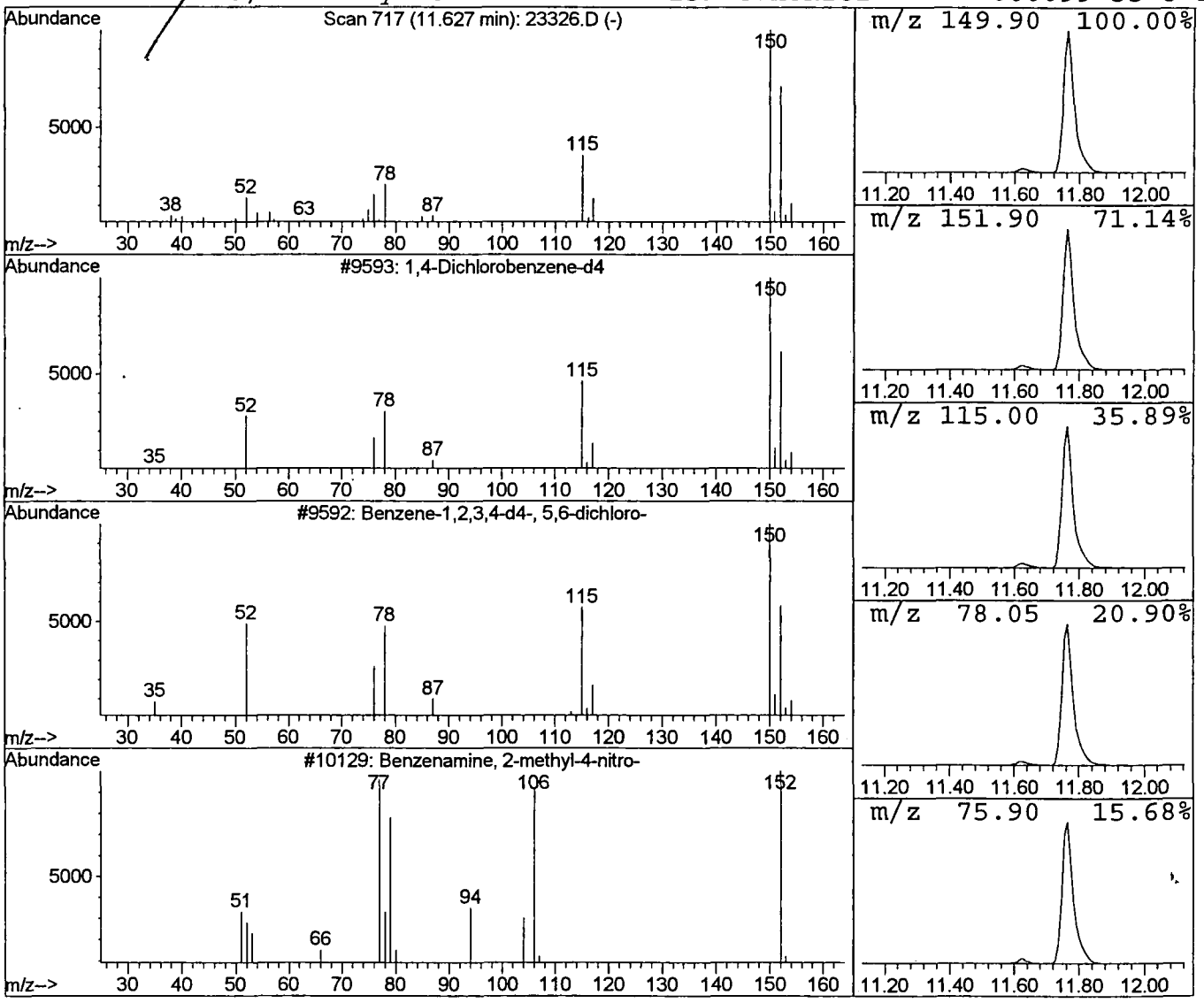
Quant Method : C:\HPCHEM\1\METHODS\NP091101.M (RTE Integrator)
Title : 209 625/TCLP calibration table 7/16/01
Library : C:\DATABASE\NBS75K.L

**********

Peak Number 6 1,4-Dichlorobenzene-d4 Concentration Rank 3

| R.T. | EstConc | Area | Relative to ISTD. | R.T. |
|---|---|---|---|---|
| 11.63 | 0.55 ug/l | 50465 | 1,4-Dichlorobenzene-d4 | 11.76 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | 1,4-Dichlorobenzene-d4 | 150 | C6Cl2D4 | 003855-82-1 | 80 |
| 2 | Benzene-1,2,3,4-d4-, 5,6-dichloro- | 150 | C6Cl2D4 | 002199-69-1 | 33 |
| 3 | Benzenamine, 2-methyl-4-nitro- | 152 | C7H8N2O2 | 000099-52-5 | 12 |
| 4 | Benzenamine, 2-methyl-5-nitro- | 152 | C7H8N2O2 | 000099-55-8 | 10 |



# Tentatively Identified Compound (LSC) summary

Operator ID: 209 GALOS Date Acquired: 10 Oct 2001 1:04 am
Data File: C:\HPCHEM\1\DATA\OCT0901\23326.D
Name: gc/ms unknown
Misc:
Method: C:\HPCHEM\1\METHODS\NP091101.M (RTE Integrator)
Title: 209 625/TCLP calibration table 7/16/01
Library Searched: C:\DATABASE\NBS75K.L

| TIC Top Hit name | RT | EstConc | Units | Area | IntStd | ISRT | ISArea | ISConc |
|---|---|---|---|---|---|---|---|---|
| 3-Hexanone | 6.38 | 0.4 | ug/l | 41168 | ISTD01 | 11.76 | 1830350 | 20.0 |
| 2-Hexanone | 6.49 | 0.9 | ug/l | 82182 | ISTD01 | 11.76 | 1830350 | 20.0 |
| 2-Pentanol, 4-methyl | 6.73 | 0.6 | ug/l | 50824 | ISTD01 | 11.76 | 1830350 | 20.0 |
| 1-Propanamine, 2-met | 10.21 | 0.4 | ug/l | 40208 | ISTD01 | 11.76 | 1830350 | 20.0 |
| Oxirane, 2-methyl-3- | 10.42 | 0.5 | ug/l | 46674 | ISTD01 | 11.76 | 1830350 | 20.0 |
| 1,4-Dichlorobenzene- | 11.63 | 0.6 | ug/l | 50465 | ISTD01 | 11.76 | 1830350 | 20.0 |

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