User: Galos, Marie
Westchester Dept. of Labs & Research
Date: 12/04/2001 Time: 15:12:05

Sample: AD26387
Analysis: $GCMS GCMS Scan For Unknowns
Units: ug
Started: 12/4/01 15:11 Ended: 12/4/01 15:11 Analyst: MG
Page: 1

	Component Name	Viol	Result
1	Other unknown compounds		see comment

Comments for Sample AD26387 - Analysis $GCMS

Westchester Dept. of Labs & Research
User: Galos, Marie
Date: 12-04-2001 Time: 15:12:28

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\NOV3001\26387.D Vial: 13
Acq On : 30 Nov 2001 11:08 pm Operator: GALOS
Sample : gc/ms unknown 11/3 Inst : GC/MS Ins
Misc : Multiplr: 1.00
MS Integration Params: RTEINT.P
Quant Time: Dec 3 15:24 2001 Quant Results File: NP112701.RES

Quant Method : C:\HPCHEM\1\METHODS\NP112701.M (RTE Integrator)
Title : 208 625/TCLP calibration table
Last Update : Wed Nov 28 15:33:44 2001
Response via : Initial Calibration
DataAcq Meth : NP112701

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	11.89	152	1591272	20.00	ng/uL	-0.05
19) Naphthalene-d8	15.50	136	6246516	20.00	ng/uL	-0.05
31) Acenaphthene-d10	20.78	164	2812440	20.00	ng/uL	-0.05
49) Phenanthrene-d10	25.19	188	4087037	20.00	ng/uL	-0.06
59) Chrysene-d12	33.19	240	2582760	20.00	ng/uL	-0.06
68) Perylene-d12	37.28	264	1824567	20.00	ng/uL	-0.06
System Monitoring Compounds						
4) 2-Fluorophenol	0.00	112	0	0.00	ng/uL	
Spiked Amount 75.000	Range 18 - 42		Recovery = 0.00%#			
5) Phenol-d5	11.18	99	2175	0.02	ng/uL	0.05
Spiked Amount 75.000	Range 19 - 32		Recovery = 0.03%#			
6) 2-Chlorophenol-d4	11.89	132	1291	0.01	ng/uL	0.47
Spiked Amount 75.000	Range 34 - 84		Recovery = 0.01%#			
7) 1,2-Dichlorobenzene-d4	12.40	152	15705	0.23	ng/uL	-0.06
Spiked Amount 50.000	Range 26 - 73		Recovery = 0.46%#			
20) Nitrobenzene-d5	0.00	82	0	0.00	ng/uL	
Spiked Amount 50.000	Range 55 - 98		Recovery = 0.00%#			
32) 2-Fluorobiphenyl	18.92	172	5630	0.03	ng/uL	0.09
Spiked Amount 50.000	Range 42 - 78		Recovery = 0.06%#			
33) 2,4,6-Tribromophenol	0.00	330	0	0.00	ng/uL	
Spiked Amount 75.000	Range 45 - 96		Recovery = 0.00%#			
62) Terphenyl-d14	0.00	244	0	0.00	ng/uL	
Spiked Amount 50.000	Range 58 - 98		Recovery = 0.00%#			
Target Compounds						Qvalue
2) Pyridine	0.00	79	0	N.D.		
3) N-nitrosodimethylamine	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	12.88	45	1425	N.D.		
16) 3&4-Methylphenol	0.00	108	0	N.D.		
17) N-Nitrosodi-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
6387.D NP112701.M Tue Dec 04 09:21:29 2001

Quantitation Report (QT Reviewed)

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Acq On : 30 Nov 2001 11:08 pm Operator: GALOS
Sample : gc/ms unknown 11/3 Inst : GC/MS Ins
Misc : Multiplr: 1.00
MS Integration Params: RTEINT.P
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Quant Method : C:\HPCHEM\1\METHODS\NP112701.M (RTE Integrator)
Title : 208 625/TCLP calibration table
Last Update : Wed Nov 28 15:33:44 2001
Response via : Initial Calibration
DataAcq Meth : NP112701

	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-Chlorophenylphenylether	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	169	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	25.19	178	277	N.D.	
56)	Anthracene	25.19	178	277	N.D.	
57)	Di-n-butylphthalate	27.18	149	7135	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	33.19	228	5648	N.D.	
66)	Bis(2-ethylhexyl)phthalate	33.45	149	14078	N.D.	
67)	Chrysene	33.19	228	5648	N.D.	
69)	Di-n-Octylphthalate	0.00	149	0	N.D.	
70)	Benzo(b)fluoranthene	0.00	252	0	N.D.	

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Misc : Multiplr: 1.00
MS Integration Params: RTEINT.P
Quant Time: Dec 3 15:24 2001 Quant Results File: NP112701.RES

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Title : 208 625/TCLP calibration table
Last Update : Wed Nov 28 15:33:44 2001
Response via : Initial Calibration
DataAcq Meth : NP112701

Compound	R.T.	QIon	Response	Conc Unit	Qvalue
71) Benzo(k)fluoranthene	0.00	252	0	N.D.	
72) Benzo(a)pyrene	37.28	252	5829	N.D.	
73) Indeno(1,2,3-cd)pyrene	0.00	276	0	N.D.	
74) Dibenz(a,h)anthracene	0.00	278	0	N.D.	
75) Benzo(g,h,i)perylene	0.00	276	0	N.D.	

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Quantitation Report

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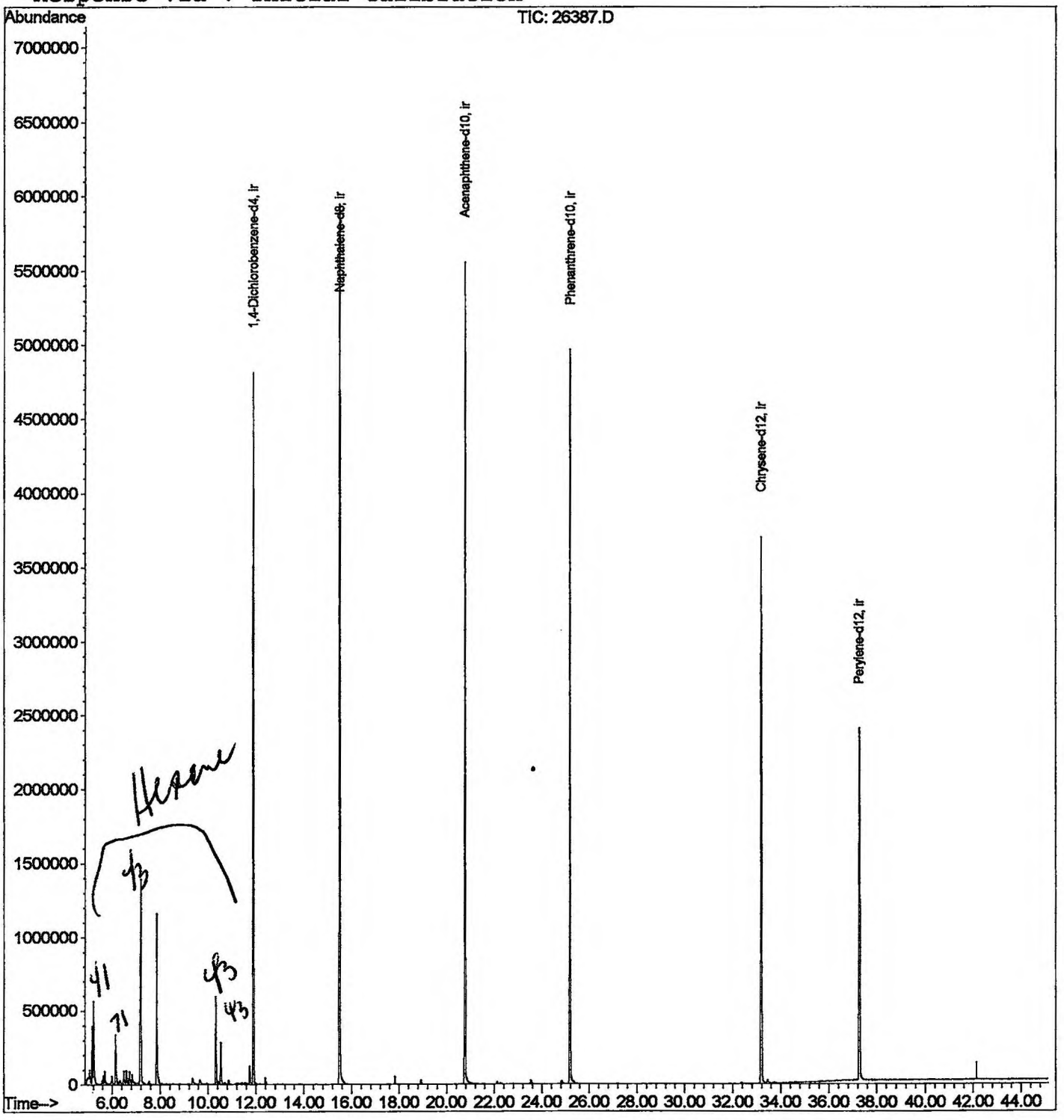
Method : C:\HPCHEM\1\METHODS\NP112701.M (RTE Integrator)
Title : 208 625/TCLP calibration table
Last Update : Wed Nov 28 15:33:44 2001
Response via : Initial Calibration



LSC Area Percent Report

Data File : C:\HPCHEM\1\DATA\NOV3001\26387.D Vial: 13
Acq On : 30 Nov 2001 11:08 pm Operator: GALOS
Sample : gc/ms unknown 11/3 Inst : GC/MS Ins
Misc : Multiplr: 1.00
MS Integration Params: LSCINT.P

Method : C:\HPCHEM\1\METHODS\NP112701.M (RTE Integrator)
Title : 208 625/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

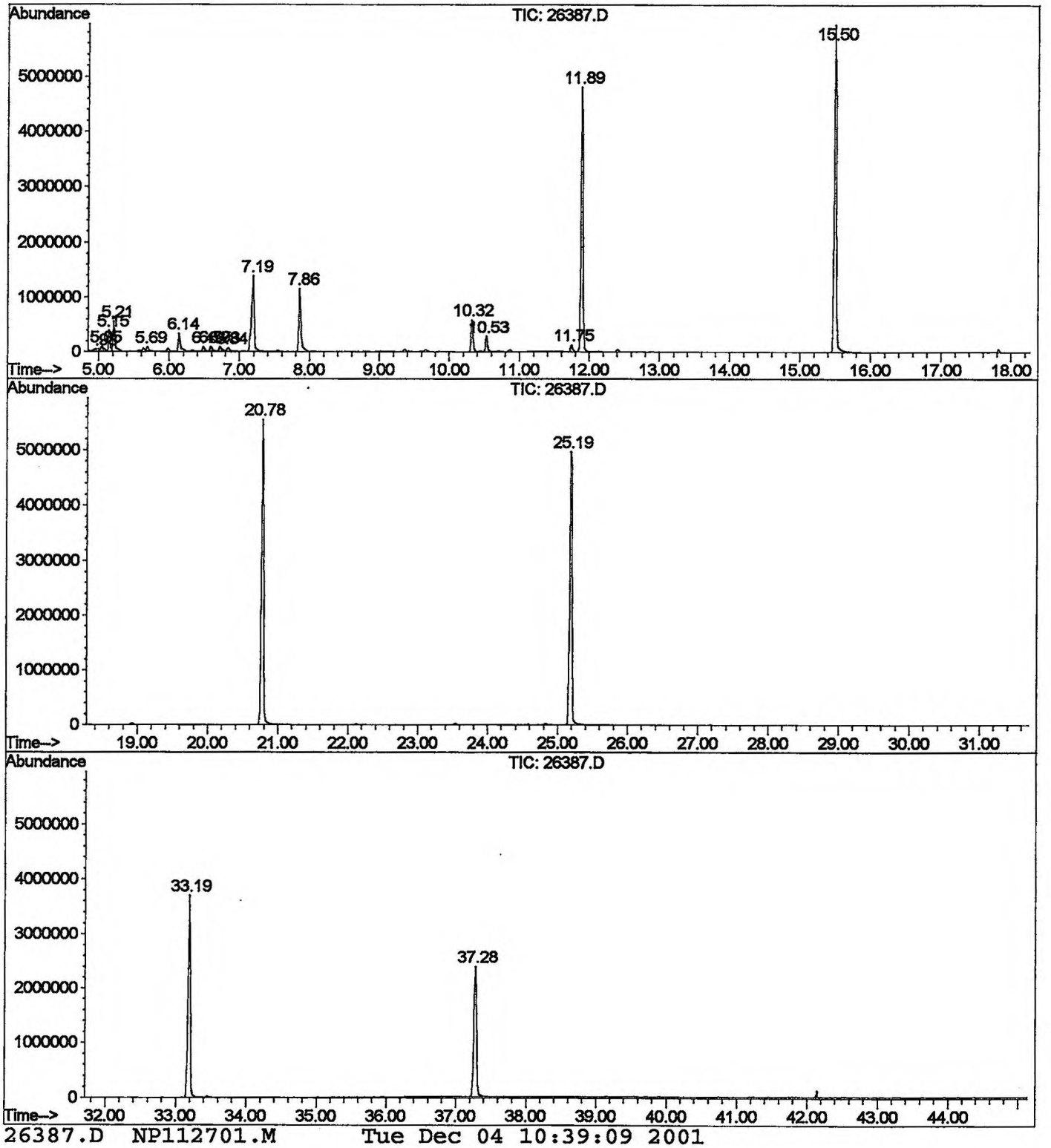
peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	peak area	peak % max.	% of total
1	4.985	6	16	19	rBV	35475	182471	1.45%	0.254%
2	5.049	19	23	31	rVV2	91539	265236	2.10%	0.369%
3	5.150	31	34	38	rVV	390953	655127	5.19%	0.911%
4	5.214	38	41	55	rVB2	561339	1110992	8.81%	1.545%
5	5.691	90	93	97	rVB	80727	138528	1.10%	0.193%
6	6.140	138	142	156	rBV	337534	719464	5.71%	1.000%
7	6.489	176	180	188	rBV	84168	167657	1.33%	0.233%
8	6.590	188	191	201	rVB	88859	190728	1.51%	0.265%
9	6.727	201	206	214	rBV2	82581	202981	1.61%	0.282%
10	6.837	214	218	223	rBV2	61325	135825	1.08%	0.189%
11	7.195	248	257	269	rBV	1386107	3464864	27.48%	4.818%
12	7.855	325	329	342	rBV	1157993	2330153	18.48%	3.240%
13	10.322	593	598	609	rBV	593254	1186015	9.40%	1.649%
14	10.533	617	621	628	rBV	282214	529313	4.20%	0.736%
15	11.752	749	754	761	rBV	125751	256985	2.04%	0.357%
16	11.890	764	769	791	rVB	4815047	9682702	76.78%	13.464%
17	15.503	1156	1163	1182	rBV	5948571	12610833	100.00%	17.535%
18	20.776	1731	1738	1759	rBV	5558550	12174920	96.54%	16.929%
19	25.187	2212	2219	2231	rBV2	4968904	11112316	88.12%	15.452%
20	33.192	3085	3092	3110	rBV2	3700369	8230567	65.27%	11.445%
21	37.282	3530	3538	3550	rBV2	2382362	6569387	52.09%	9.135%

Sum of corrected areas: 71917064

26387.D NP112701.M Tue Dec 04 10:39:08 2001

LSC Report - Integrated Chromatogram

File : C:\HPCHEM\1\DATA\NOV3001\26387.D
Operator : GALOS
Acquired : 30 Nov 2001 11:08 pm using AcqMethod NP112701
Instrument : GC/MS Ins
Sample Name: gc/ms unknown 11/3
Misc Info :
Vial Number: 13
Quant File :NP112701.RES (RTE Integrator)



26387.D NP112701.M Tue Dec 04 10:39:09 2001

Library Search Compound Report

Data File : C:\HPCHEM\1\DATA\NOV3001\26387.D
Acq On : 30 Nov 2001 11:08 pm
Sample : gc/ms unknown 11/3
Misc :
MS Integration Params: LSCINT.P

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

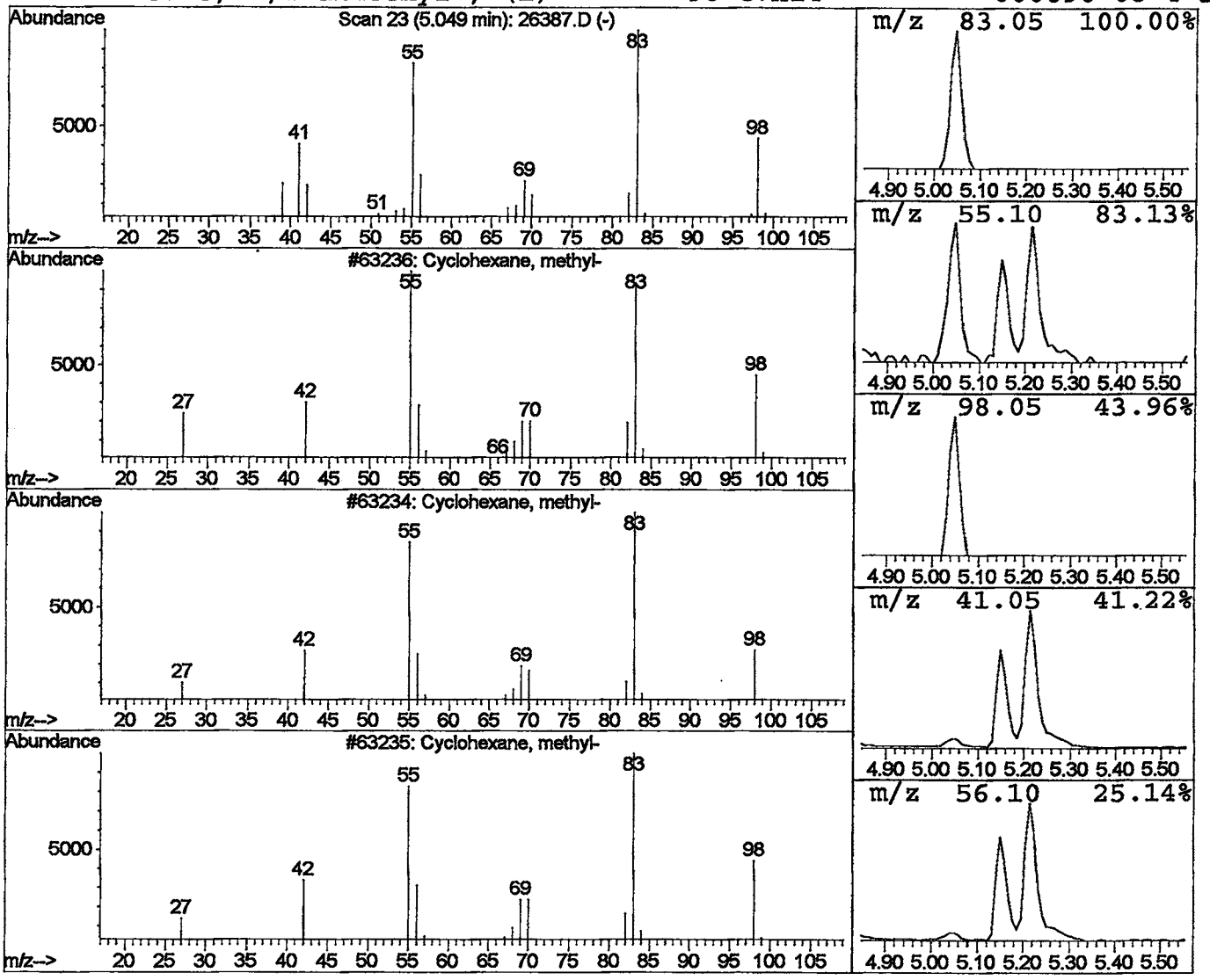
Quant Method : C:\HPCHEM\1\METHODS\NP112701.M (RTE Integrator)
Title : 208 625/TCLP calibration table
Library : C:\DATABASE\NBS75K.L

****

Peak Number 1 Cyclohexane, methyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.05	0.55 ng/uL	265236	1,4-Dichlorobenzene-d4	11.89

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclohexane, methyl-	98	C7H14	000108-87-2	81
2	Cyclohexane, methyl-	98	C7H14	000108-87-2	72
3	Cyclohexane, methyl-	98	C7H14	000108-87-2	59
4	2-Pentene, 4,4-dimethyl-, (E)-	98	C7H14	000690-08-4	53



Library Search Compound Report

Data File : C:\HPCHEM\1\DATA\NOV3001\26387.D
Acq On : 30 Nov 2001 11:08 pm
Sample : gc/ms unknown 11/3
Misc :
MS Integration Params: LSCINT.P

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

Quant Method : C:\HPCHEM\1\METHODS\NP112701.M (RTE Integrator)
Title : 208 625/TCLP calibration table
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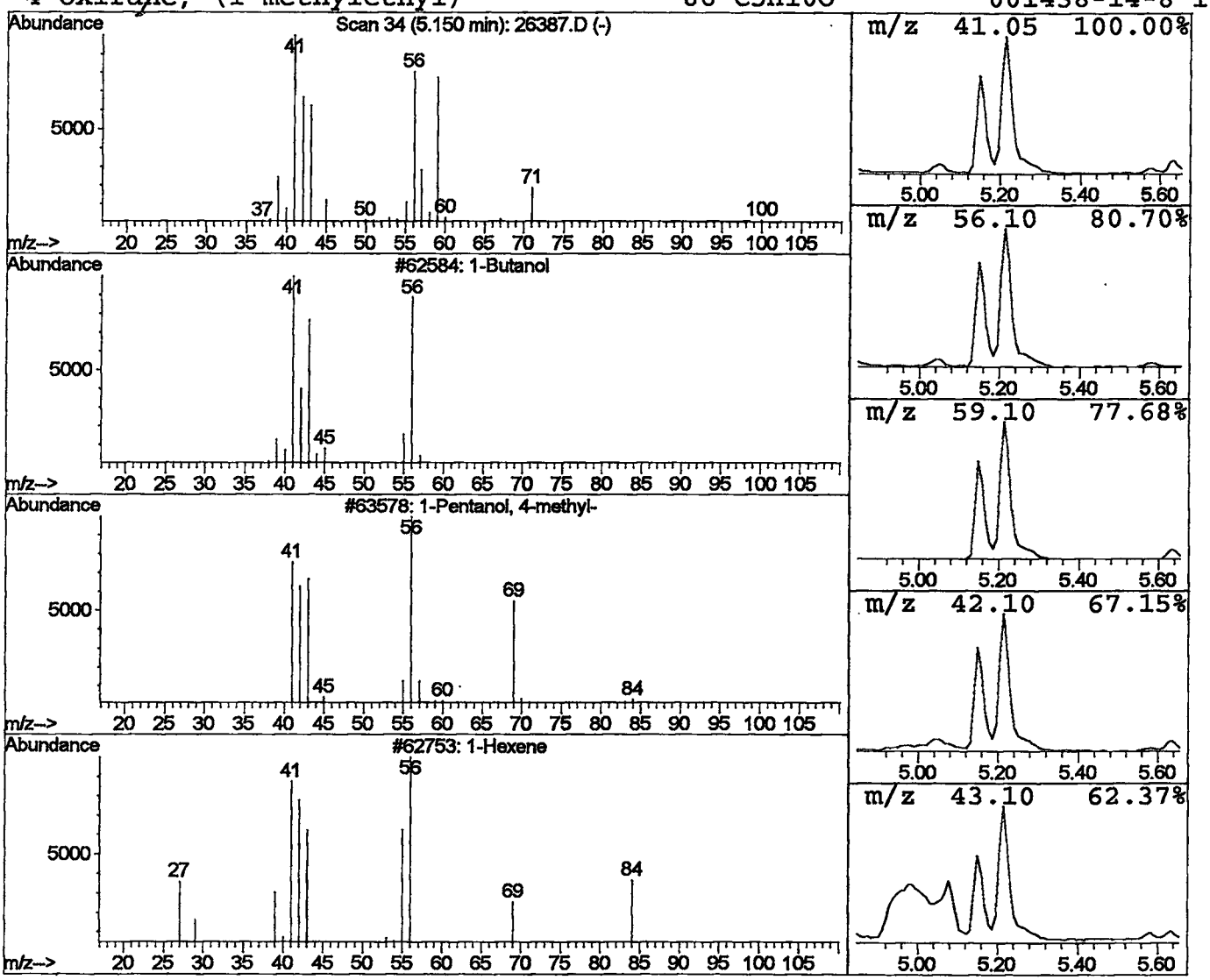
****

Peak Number 2 1-Butanol Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.15	1.35 ng/uL	655127	1,4-Dichlorobenzene-d4	11.89

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Butanol	74	C4H10O	000071-36-3	32
2	1-Pentanol, 4-methyl-	102	C6H14O	000626-89-1	25
3	1-Hexene	84	C6H12	000592-41-6	25
4	Oxirane, (1-methylethyl)-	86	C5H10O	001438-14-8	16



Library Search Compound Report

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Acq On : 30 Nov 2001 11:08 pm
Sample : gc/ms unknown 11/3
Misc :
MS Integration Params: LSCINT.P

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

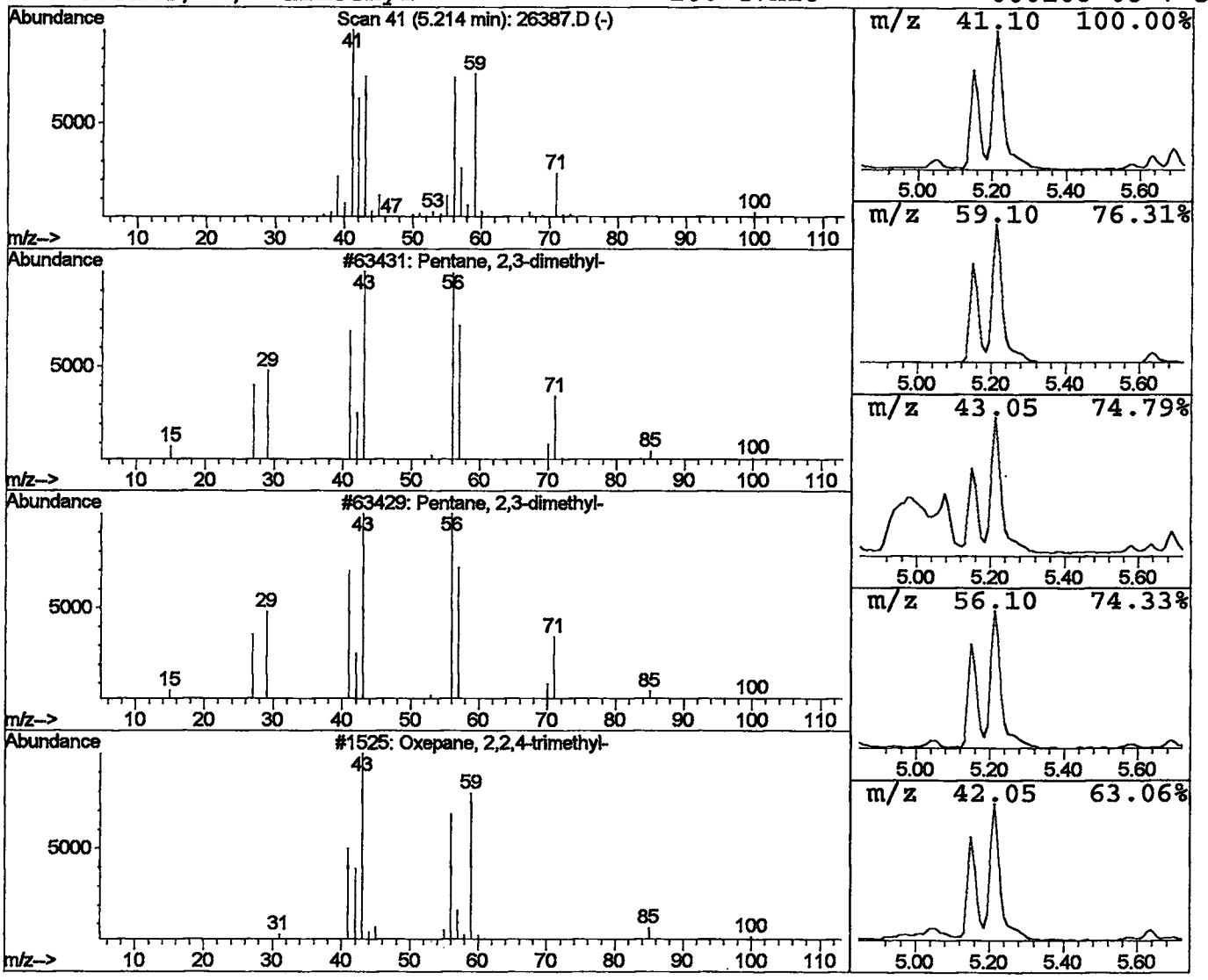
Quant Method : C:\HPCHEM\1\METHODS\NP112701.M (RTE Integrator)
Title : 208 625/TCLP calibration table
Library : C:\DATABASE\NBS75K.L

****

Peak Number 3 Pentane, 2,3-dimethyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.21	2.29 ng/uL	1110990	1,4-Dichlorobenzene-d4	11.89

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Pentane, 2,3-dimethyl-	100	C7H16	000565-59-3	38
2	Pentane, 2,3-dimethyl-	100	C7H16	000565-59-3	38
3	Oxepane, 2,2,4-trimethyl-	100	C6H12O	023120-44-7	36
4	Pentane, 2,4-dimethyl-	100	C7H16	000108-08-7	35



Library Search Compound Report

Data File : C:\HPCHEM\1\DATA\NOV3001\26387.D
Acq On : 30 Nov 2001 11:08 pm
Sample : gc/ms unknown 11/3
Misc :
MS Integration Params: LSCINT.P

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

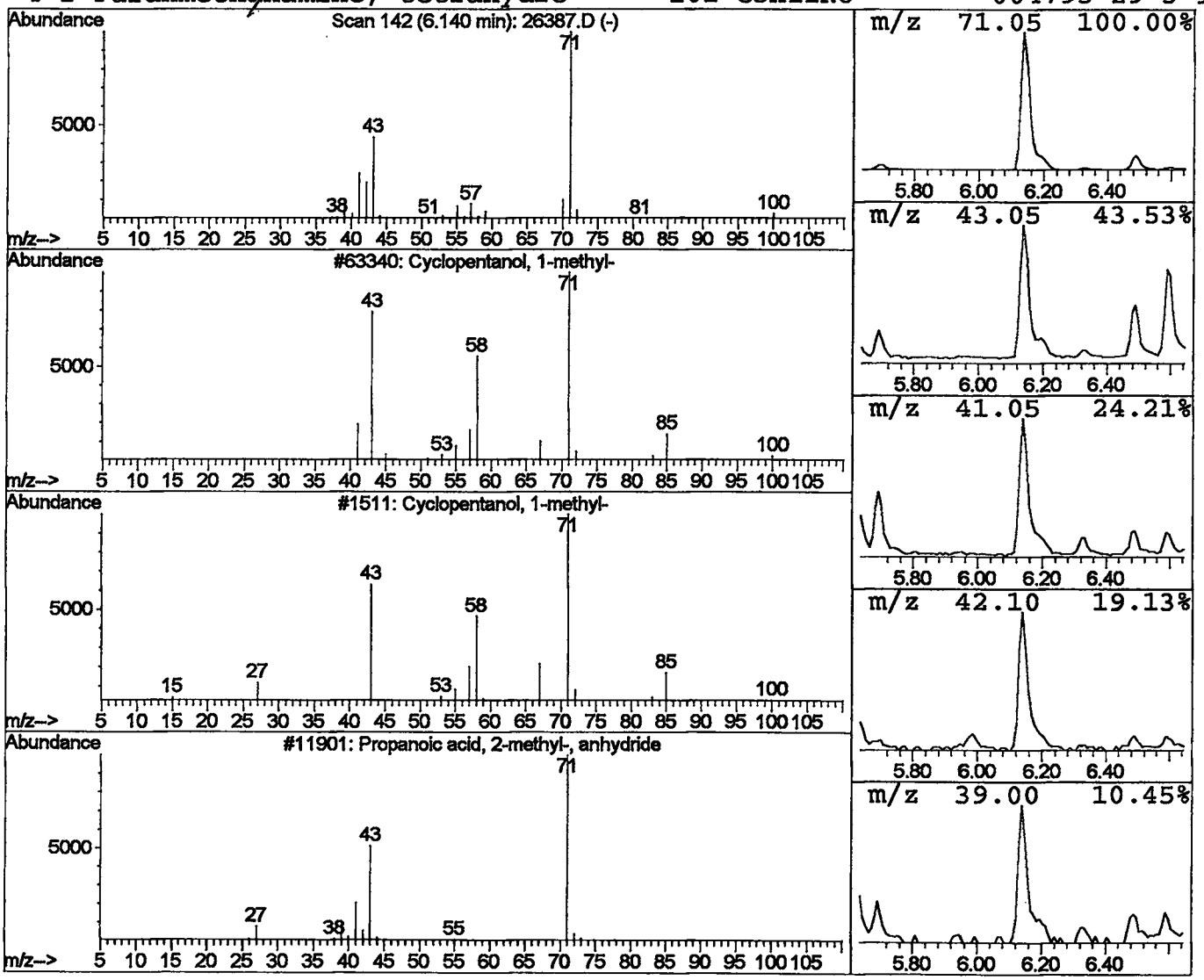
Quant Method : C:\HPCHEM\1\METHODS\NP112701.M (RTE Integrator)
Title : 208 625/TCLP calibration table
Library : C:\DATABASE\NBS75K.L

****

Peak Number 4 Cyclopentanol, 1-methyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.14	1.49 ng/uL	719464	1,4-Dichlorobenzene-d4	11.89

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclopentanol, 1-methyl-	100	C6H12O	001462-03-9	28
2	Cyclopentanol, 1-methyl-	100	C6H12O	001462-03-9	9
3	Propanoic acid, 2-methyl-, anhydrid	158	C8H14O3	000097-72-3	9
4	2-Furanmethanamine, tetrahydro-	101	C5H11NO	004795-29-3	9



Library Search Compound Report

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Operator: GALOS
Inst : GC/MS Ins
Multiplr: 1.00

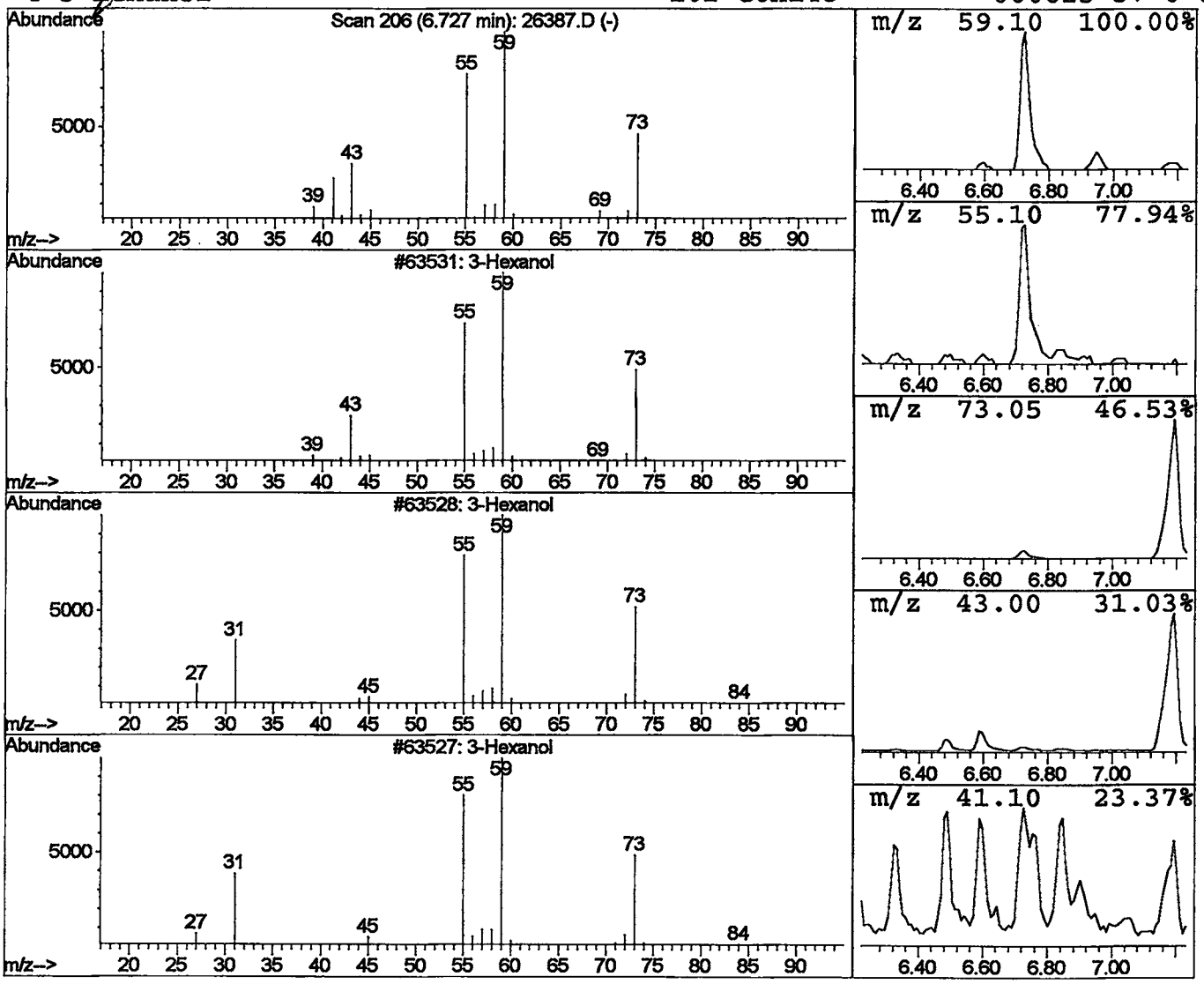
Quant Method : C:\HPCHEM\1\METHODS\NP112701.M (RTE Integrator)
Title : 208 625/TCLP calibration table
Library : C:\DATABASE\NBS75K.L

****

Peak Number 5 3-Hexanol — Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.73	0.42 ng/uL	202981	1,4-Dichlorobenzene-d4	11.89

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	3-Hexanol	102	C6H14O	000623-37-0	83
2	3-Hexanol	102	C6H14O	000623-37-0	83
3	3-Hexanol	102	C6H14O	000623-37-0	78
4	3-Hexanol	102	C6H14O	000623-37-0	64



Library Search Compound Report

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Sample : gc/ms unknown 11/3
Misc :
MS Integration Params: LSCINT.P

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

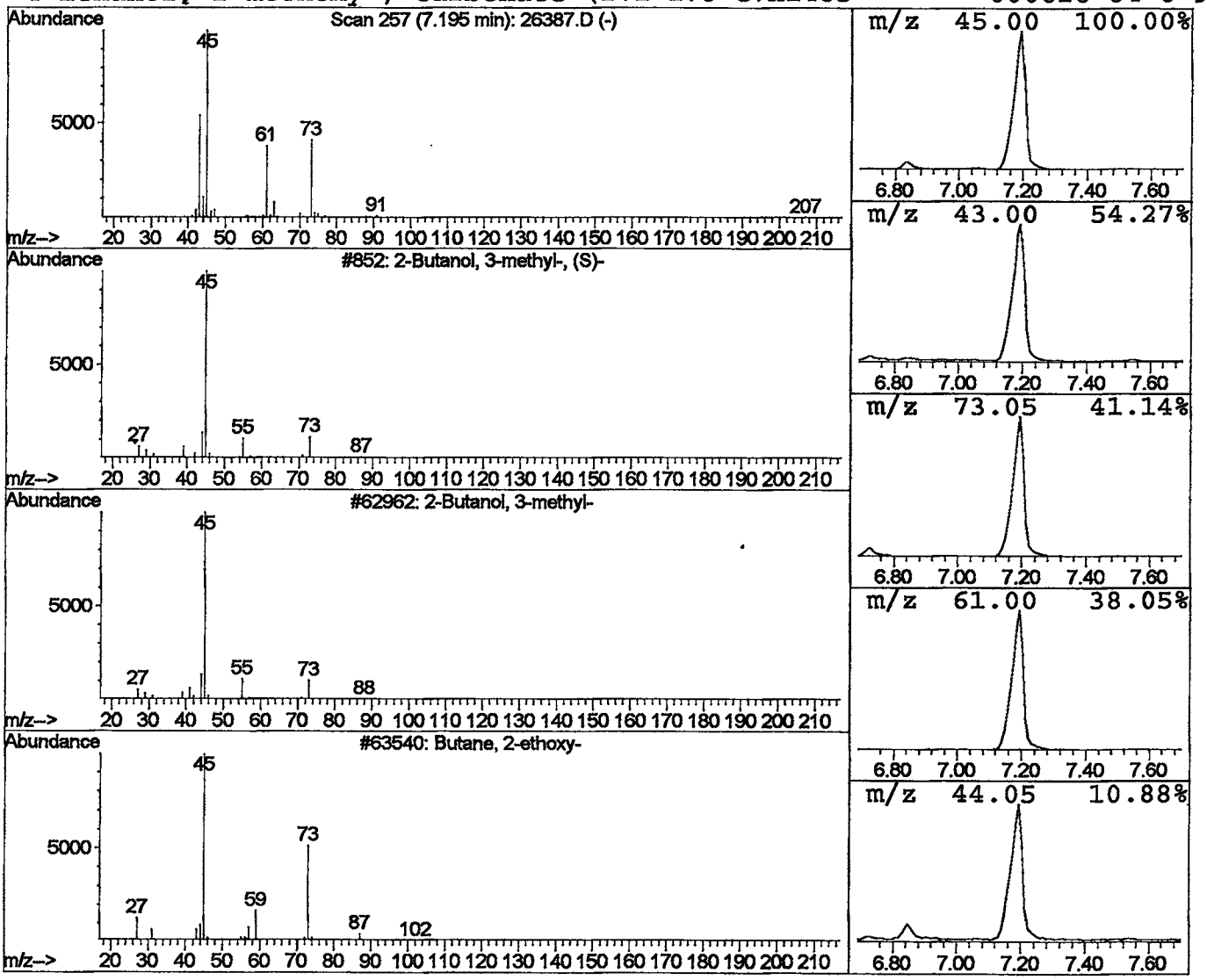
Quant Method : C:\HPCHEM\1\METHODS\NP112701.M (RTE Integrator)
Title : 208 625/TCLP calibration table
Library : C:\DATABASE\NBS75K.L

****

Peak Number 6 2-Butanol, 3-methyl-, (S)- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.19	7.16 ng/uL	3464860	1,4-Dichlorobenzene-d4	11.89

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Butanol, 3-methyl-, (S)-	88	C5H12O	001517-66-4	38
2	2-Butanol, 3-methyl-	88	C5H12O	000598-75-4	9
3	Butane, 2-ethoxy-	102	C6H14O	002679-87-0	9
4	Ethanol, 2-methoxy-, carbonate (2:1	178	C7H14O5	000626-84-6	9



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

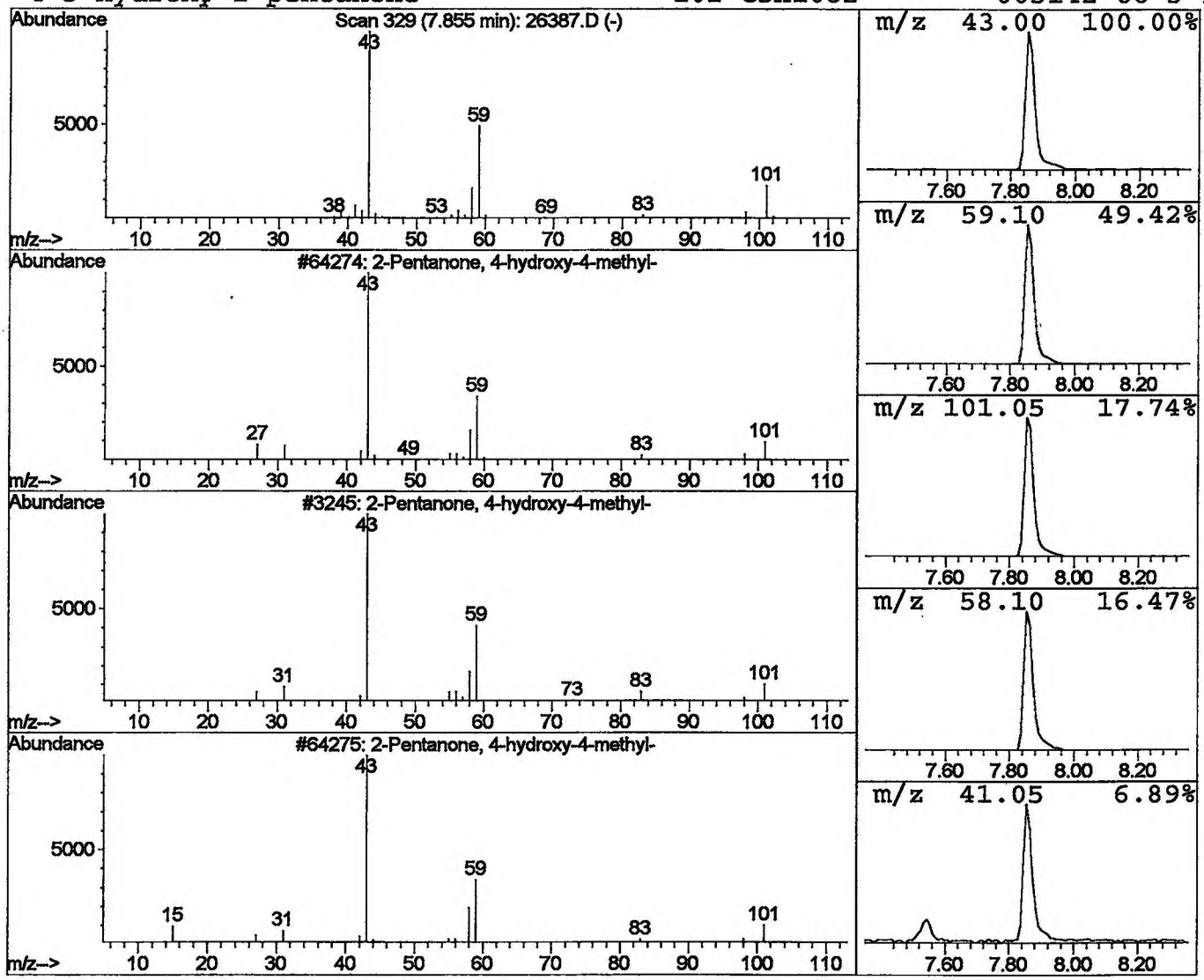
Quant Method : C:\HPCHEM\1\METHODS\NP112701.M (RTE Integrator)
Title : 208 625/TCLP calibration table
Library : C:\DATABASE\NBS75K.L

****

Peak Number 7 2-Pentanone, 4-hydroxy-4-methy Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.86	4.81 ng/uL	2330150	1,4-Dichlorobenzene-d4	11.89

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	59
2	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	43
3	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	25
4	3-Hydroxy-2-pentanone	102	C5H10O2	003142-66-3	9



Library Search Compound Report

Data File : C:\HPCHEM\1\DATA\NOV3001\26387.D
Acq On : 30 Nov 2001 11:08 pm
Sample : gc/ms unknown 11/3
Misc :
MS Integration Params: LSCINT.P

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

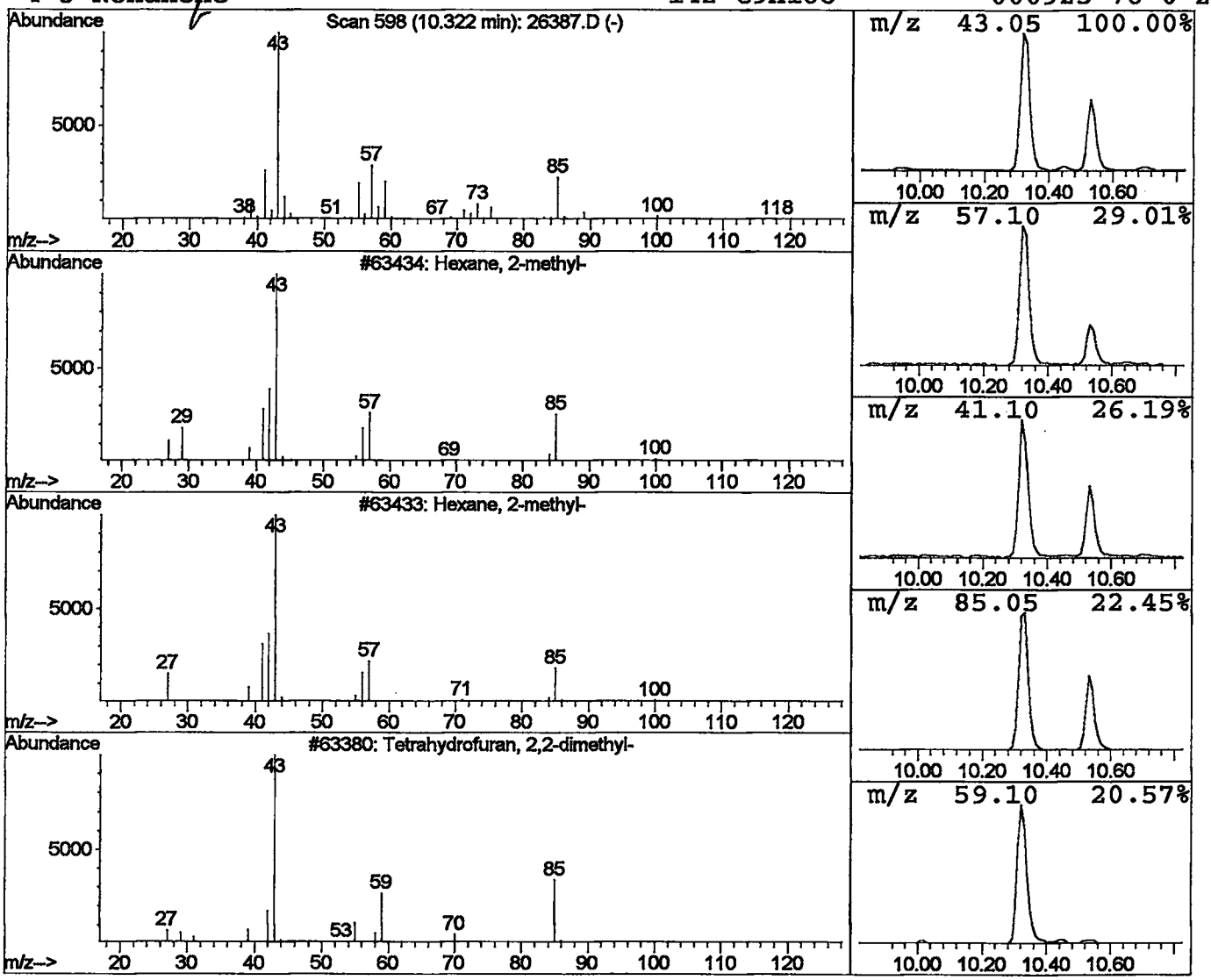
Quant Method : C:\HPCHEM\1\METHODS\NP112701.M (RTE Integrator)
Title : 208 625/TCLP calibration table
Library : C:\DATABASE\NBS75K.L

****

Peak Number 8 Hexane, 2-methyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.32	2.45 ng/uL	1186020	1,4-Dichlorobenzene-d4	11.89

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Hexane, 2-methyl-	100	C7H16	000591-76-4	35
2	Hexane, 2-methyl-	100	C7H16	000591-76-4	32
3	Tetrahydrofuran, 2,2-dimethyl-	100	C6H12O	001003-17-4	28
4	3-Nonanone	142	C9H18O	000925-78-0	25



Library Search Compound Report

Data File : C:\HPCHEM\1\DATA\NOV3001\26387.D
Acq On : 30 Nov 2001 11:08 pm
Sample : gc/ms unknown 11/3
Misc :
MS Integration Params: LSCINT.P

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

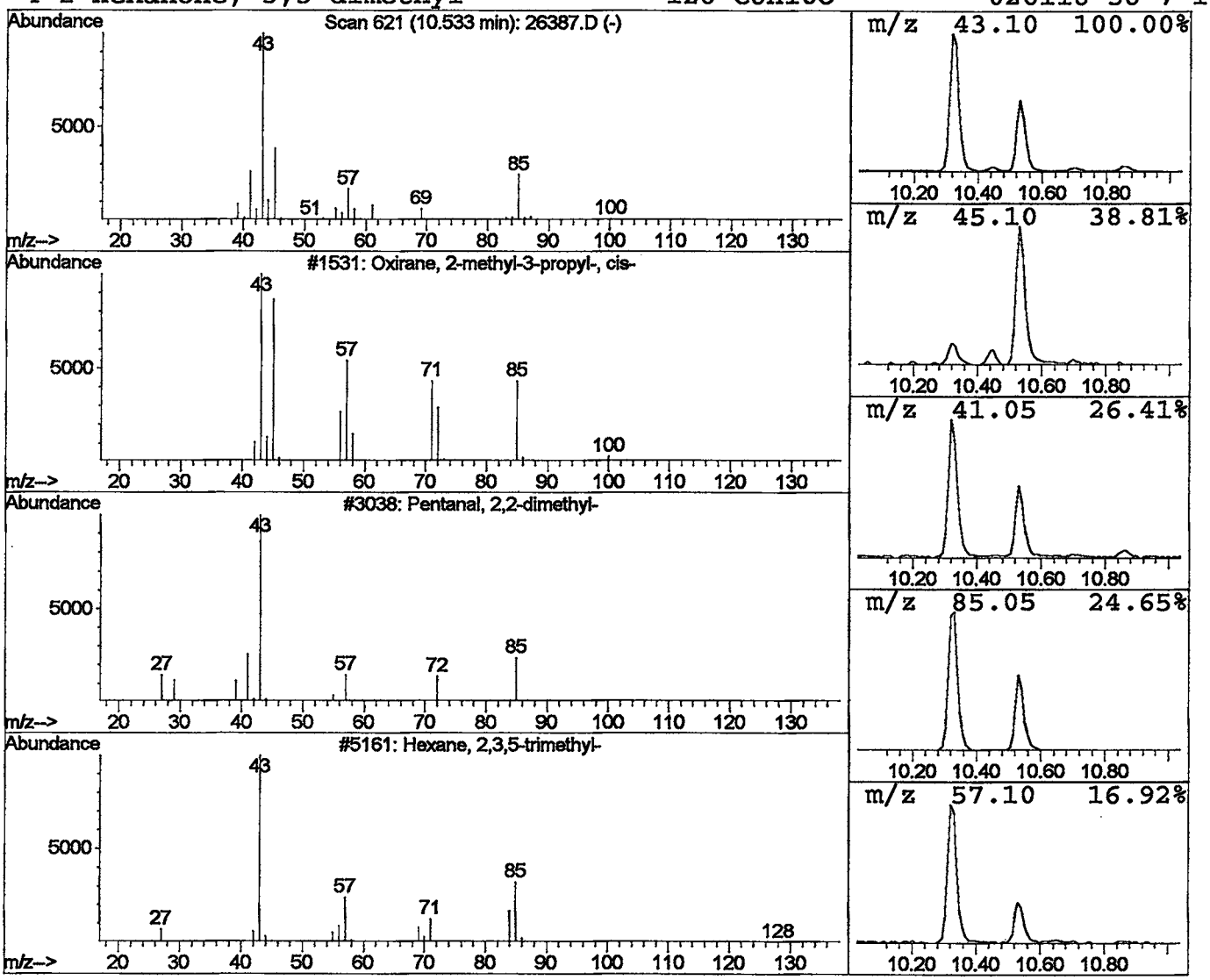
Quant Method : C:\HPCHEM\1\METHODS\NP112701.M (RTE Integrator)
Title : 208 625/TCLP calibration table
Library : C:\DATABASE\NBS75K.L

****

Peak Number 9 Oxirane, 2-methyl-3-propyl-, c Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.53	1.09 ng/uL	529313	1,4-Dichlorobenzene-d4	11.89

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Oxirane, 2-methyl-3-propyl-, cis-	100	C6H12O	006124-90-9	50
2	Pentanal, 2,2-dimethyl-	114	C7H14O	014250-88-5	25
3	Hexane, 2,3,5-trimethyl-	128	C9H20	001069-53-0	17
4	2-Hexanone, 3,3-dimethyl-	128	C8H16O	026118-38-7	17



Library Search Compound Report

Data File : C:\HPCHEM\1\DATA\NOV3001\26387.D Vial: 13
Acq On : 30 Nov 2001 11:08 pm Operator: GALOS
Sample : gc/ms unknown 11/3 Inst : GC/MS Ins
Misc : Multiplr: 1.00
MS Integration Params: LSCINT.P

Quant Method : C:\HPCHEM\1\METHODS\NP112701.M (RTE Integrator)
Title : 208 625/TCLP calibration table
Library : C:\DATABASE\NBS75K.L

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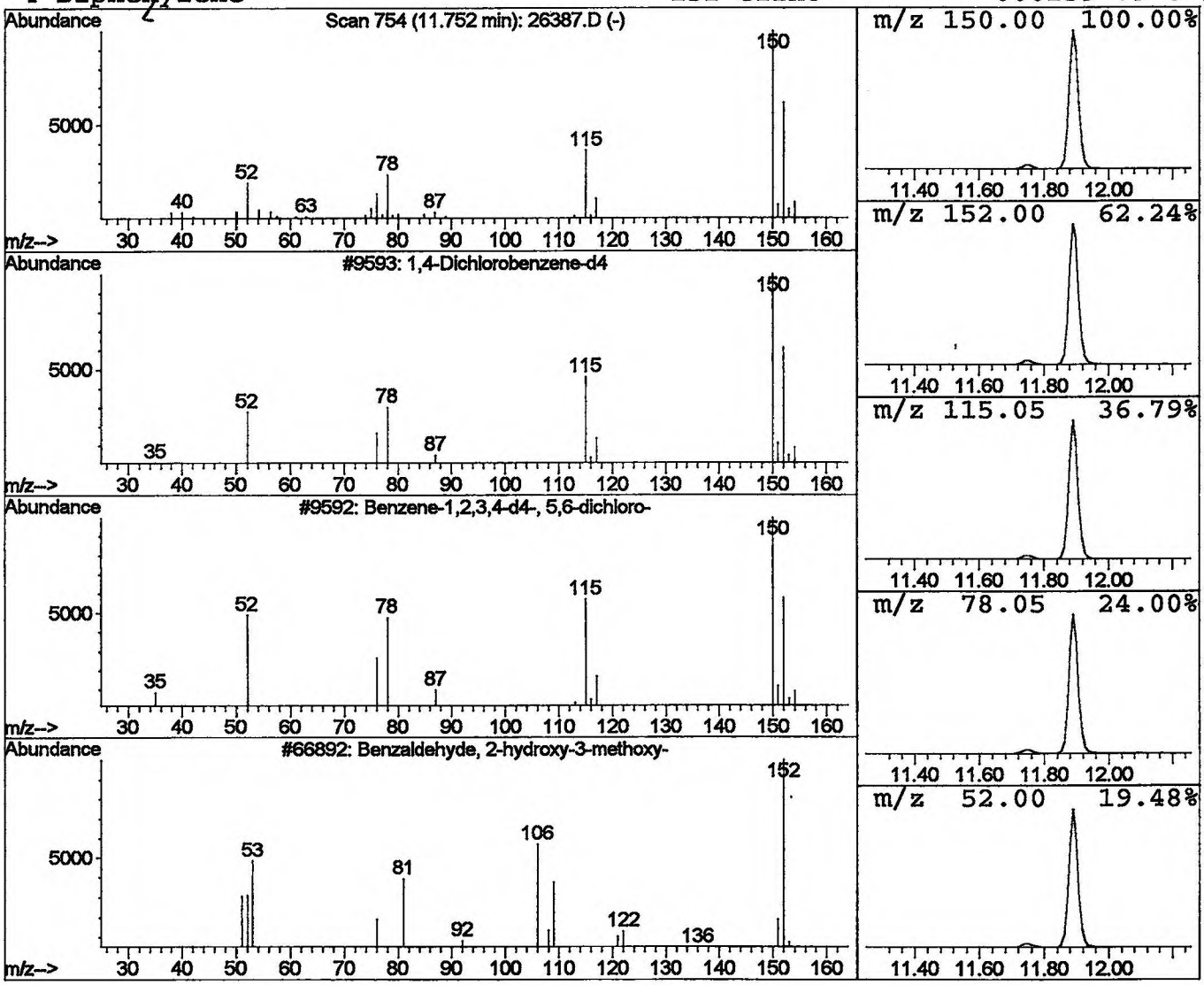
Peak Number 10 1,4-Dichlorobenzene-d4 Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.75	0.53 ng/uL	256985	1,4-Dichlorobenzene-d4	11.89

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,4-Dichlorobenzene-d4	150	C6Cl2D4	003855-82-1	59
2	Benzene-1,2,3,4-d4-, 5,6-dichloro-	150	C6Cl2D4	002199-69-1	45
3	Benzaldehyde, 2-hydroxy-3-methoxy-	152	C8H8O3	000148-53-8	35
4	Biphenylene	152	C12H8	000259-79-0	9



Tentatively Identified Compound (LSC) summary

Operator ID: GALOS Date Acquired: 30 Nov 2001 11:08 pm
Data File: C:\HPCHEM\1\DATA\NOV3001\26387.D
Name: gc/ms unknown 11/3
Misc:
Method: C:\HPCHEM\1\METHODS\NP112701.M (RTE Integrator)
Title: 208 625/TCLP calibration table
Library Searched: C:\DATABASE\NBS75K.L

TIC Top Hit name	RT	EstConc	Units	Area	IntStd	ISRT	ISArea	ISConc
Cyclohexane, methyl-	5.05	0.5	ng/uL	265236	ISTD01	11.89	9682700	20.0
1-Butanol	5.15	1.4	ng/uL	655127	ISTD01	11.89	9682700	20.0
Pentane, 2,3-dimethy	5.21	2.3	ng/uL	1110990	ISTD01	11.89	9682700	20.0
Cyclopentanol, 1-met	6.14	1.5	ng/uL	719464	ISTD01	11.89	9682700	20.0
3-Hexanol	6.73	0.4	ng/uL	202981	ISTD01	11.89	9682700	20.0
2-Butanol, 3-methyl-	7.19	7.2	ng/uL	3464860	ISTD01	11.89	9682700	20.0
2-Pentanone, 4-hydro	7.86	4.8	ng/uL	2330150	ISTD01	11.89	9682700	20.0
Hexane, 2-methyl-	10.32	2.4	ng/uL	1186020	ISTD01	11.89	9682700	20.0
Oxirane, 2-methyl-3-	10.53	1.1	ng/uL	529313	ISTD01	11.89	9682700	20.0
1,4-Dichlorobenzene-	11.75	0.5	ng/uL	256985	ISTD01	11.89	9682700	20.0

26387.D NP112701.M Tue Dec 04 10:39:18 2001