User: Galos, Marie
Westchester Dept. of Labs & Research
Date: 12/04/2001 Time: 15:13:07

Sample: AD26386
Analysis: $GCMS GCMS Scan For Unknowns
Units: ug
Started: 12/4/01 15:12 Ended: 12/4/01 15:12 Analyst: MG
Page: 1

	Component Name	Viol	Result
1	Other unknown compounds		see comment

Comments for Sample AD26386 - Analysis $GCMS

Westchester Dept. of Labs & Research
User: Galos, Marie
Date: 12-04-2001 Time: 15:13:17

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\26386.D Vial: 12
Acq On : 30 Nov 2001 10:13 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	1471353	20.00	ng/uL	-0.05
19) Naphthalene-d8	15.50	136	5669924	20.00	ng/uL	-0.05
31) Acenaphthene-d10	20.78	164	2515839	20.00	ng/uL	-0.05
49) Phenanthrene-d10	25.19	188	3581996	20.00	ng/uL	-0.05
59) Chrysene-d12	33.19	240	2197031	20.00	ng/uL	-0.05
68) Perylene-d12	37.28	264	1544986	20.00	ng/uL	-0.05
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	11.24	99	276	0.00	ng/uL	0.11
Spiked Amount 75.000	Range 19 - 32		Recovery =	0.00%#		
6) 2-Chlorophenol-d4	11.88	132	1149	0.01	ng/uL	0.46
Spiked Amount 75.000	Range 34 - 84		Recovery =	0.01%#		
7) 1,2-Dichlorobenzene-d4	12.40	152	15740	0.25	ng/uL	-0.05
Spiked Amount 50.000	Range 26 - 73		Recovery =	0.50%#		
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	5256	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	12.89	45	1809	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

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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
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	15.56	128	290	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	25.20	178	380	N.D.	
56)	Anthracene	25.20	178	380	N.D.	
57)	Di-n-butylphthalate	27.18	149	4789	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	4608	N.D.	
66)	Bis(2-ethylhexyl)phthalate	33.45	149	9279	N.D.	
67)	Chrysene	33.19	228	4608	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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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	4877	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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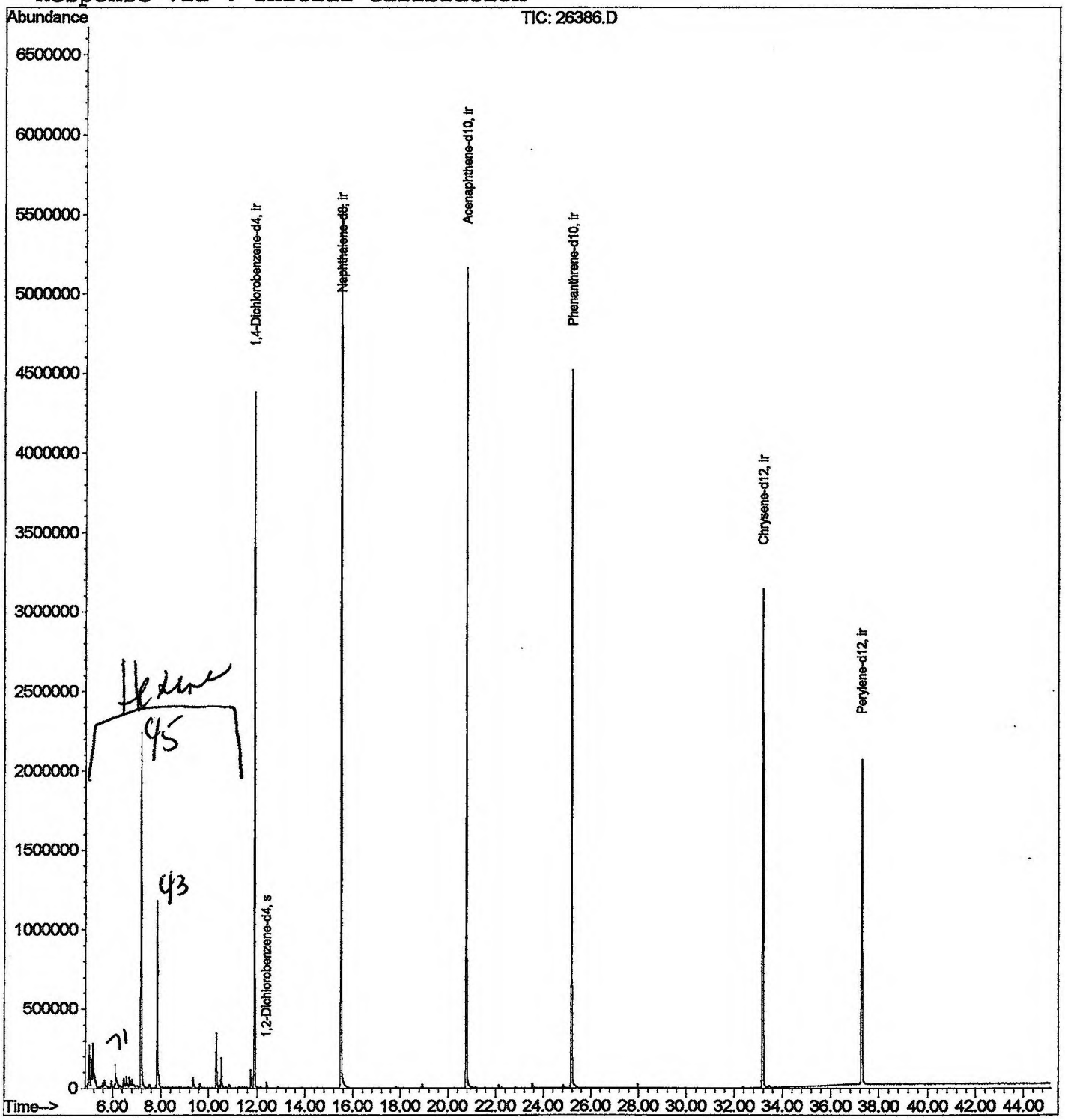
Acq On : 30 Nov 2001 10:13 pm
Sample : gc/ms unknown 11/3
Misc :
MS Integration Params: RTEINT.P
Quant Time: Dec 3 15:24 2001

Vial: 12
Operator: GALOS
Inst : GC/MS Ins
Multiplr: 1.00
Quant Results File: NP112701.RES

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\26386.D Vial: 12
Acq On : 30 Nov 2001 10:13 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

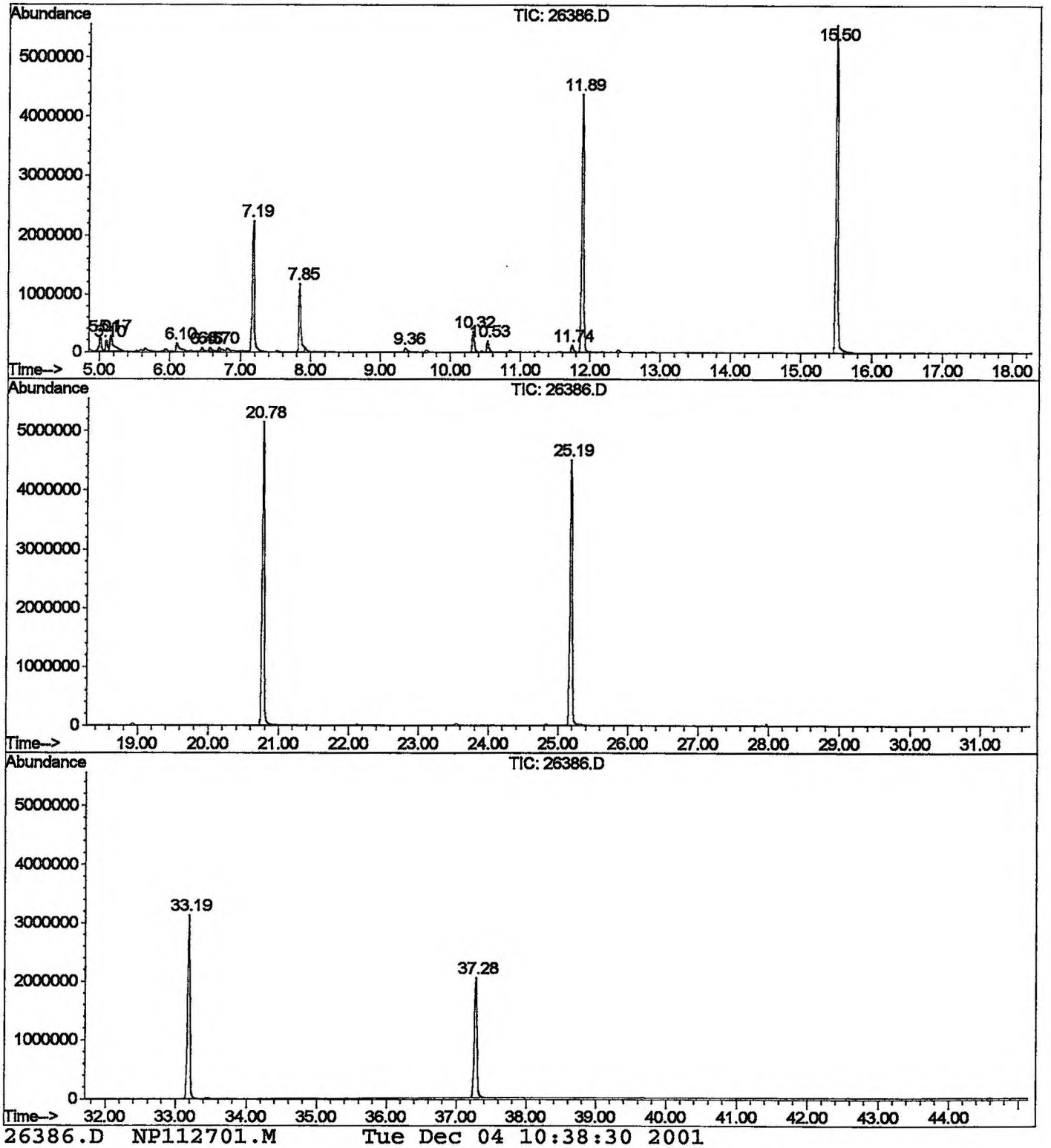
peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	peak area	peak % max.	% of total
1	5.013	10	19	25	rBV	261319	580103	5.10%	0.898%
2	5.095	25	28	32	rVV	176972	365882	3.22%	0.566%
3	5.169	32	36	55	rVB2	276900	949598	8.35%	1.469%
4	6.104	135	138	155	rBV	144358	480759	4.23%	0.744%
5	6.462	173	177	181	rBV	58687	121320	1.07%	0.188%
6	6.572	185	189	199	rVB	61280	144868	1.27%	0.224%
7	6.700	199	203	205	rBV	60055	116042	1.02%	0.180%
8	7.186	249	256	275	rVB	2237838	4480037	39.38%	6.933%
9	7.846	324	328	335	rBV	1175734	2304207	20.25%	3.566%
10	9.360	489	493	502	rBV	66566	153166	1.35%	0.237%
11	10.322	593	598	608	rBV	347691	753211	6.62%	1.166%
12	10.533	616	621	630	rBV	191233	384399	3.38%	0.595%
13	11.744	748	753	762	rVB2	111497	244101	2.15%	0.378%
14	11.891	763	769	780	rBV	4381254	8921271	78.41%	13.806%
15	15.504	1157	1163	1182	rBV	5558460	11377570	100.00%	17.607%
16	20.776	1732	1738	1757	rBV	5160487	10908346	95.88%	16.881%
17	25.187	2212	2219	2231	rBV2	4513965	9755200	85.74%	15.096%
18	33.193	3085	3092	3106	rBV2	3135759	7056194	62.02%	10.919%
19	37.282	3530	3538	3555	rBV2	2041985	5524709	48.56%	8.549%

Sum of corrected areas: 64620983

26386.D NP112701.M Tue Dec 04 10:38:29 2001

LSC Report - Integrated Chromatogram

File : C:\HPCHEM\1\DATA\NOV3001\26386.D
Operator : GALOS
Acquired : 30 Nov 2001 10:13 pm using AcqMethod NP112701
Instrument : GC/MS Ins
Sample Name: gc/ms unknown 11/3
Misc Info :
Vial Number: 12
Quant File :NP112701.RES (RTE Integrator)



26386.D NP112701.M Tue Dec 04 10:38:30 2001

Library Search Compound Report

Data File : C:\HPCHEM\1\DATA\NOV3001\26386.D
Acq On : 30 Nov 2001 10:13 pm
Sample : gc/ms unknown 11/3
Misc :
MS Integration Params: LSCINT.P

Vial: 12
Operator: GALOS
Inst : GC/MS Ins
Multiplr: 1.00

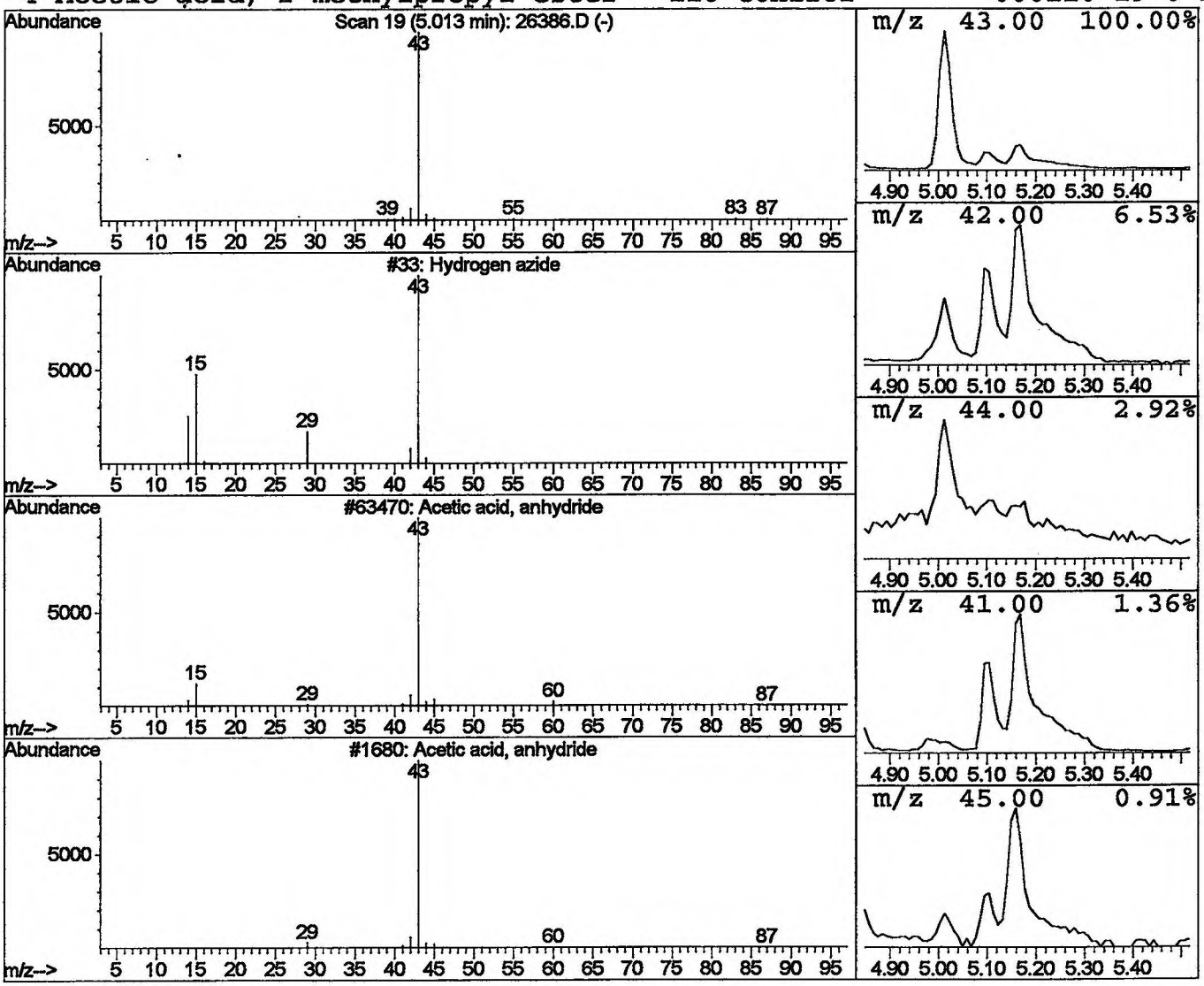
Quant Method : C:\HPCHEM\1\METHODS\NP112701.M (RTE Integrator)
Title : 208 625/TCLP calibration table
Library : C:\DATABASE\NBS75K.L

****

Peak Number 1 Hydrogen azide Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.01	1.30 ng/uL	580103	1,4-Dichlorobenzene-d4	11.89

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Hydrogen azide	43	HN3	007782-79-8	4
2	Acetic acid, anhydride	102	C4H6O3	000108-24-7	2
3	Acetic acid, anhydride	102	C4H6O3	000108-24-7	2
4	Acetic acid, 2-methylpropyl ester	116	C6H12O2	000110-19-0	2



Library Search Compound Report

Data File : C:\HPCHEM\1\DATA\NOV3001\26386.D
Acq On : 30 Nov 2001 10:13 pm
Sample : gc/ms unknown 11/3
Misc :
MS Integration Params: LSCINT.P

Vial: 12
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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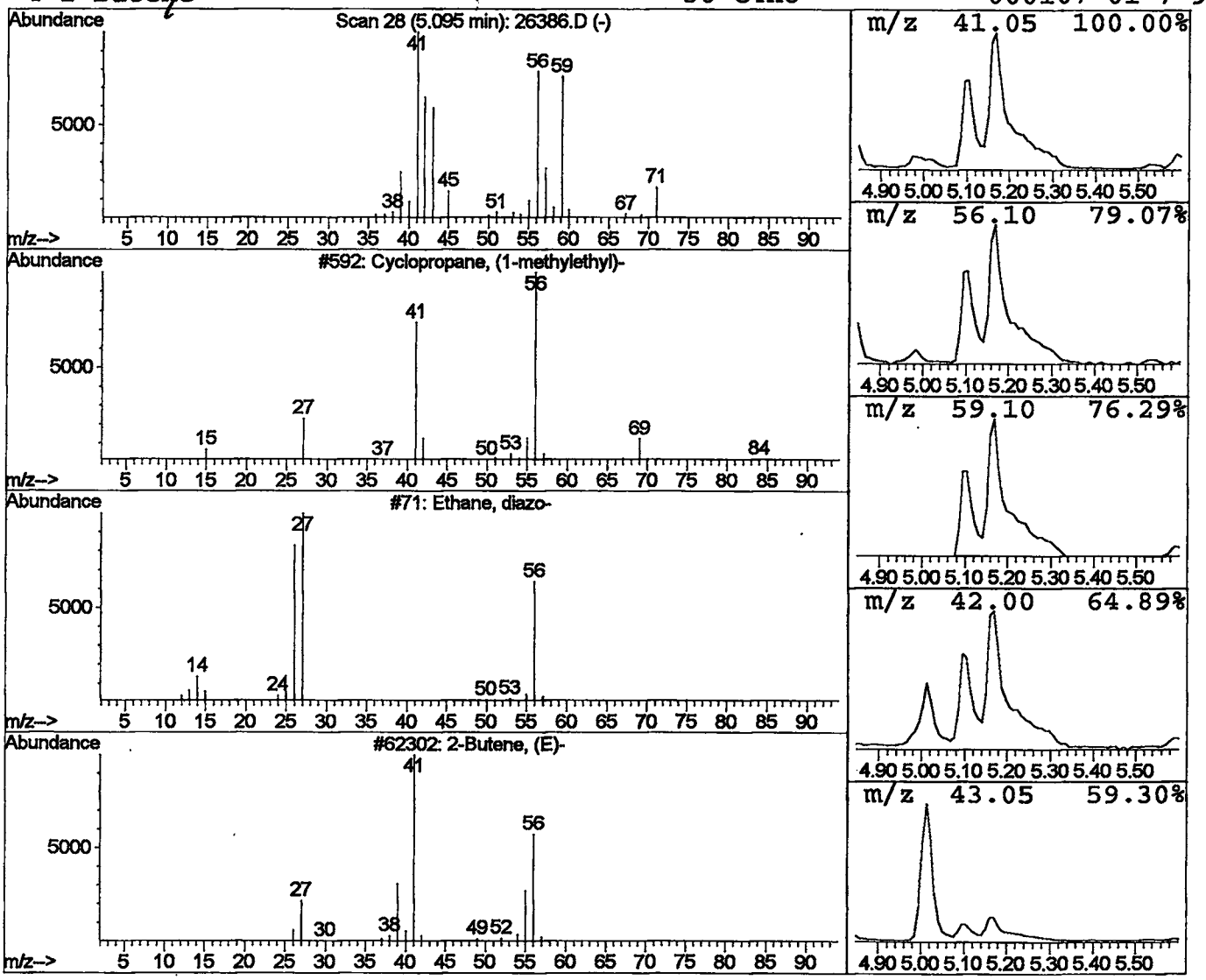
****

Peak Number 2 Cyclopropane, (1-methylethyl)- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.10	0.82 ng/uL	365882	1,4-Dichlorobenzene-d4	11.89

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclopropane, (1-methylethyl)-	84	C6H12	003638-35-5	10
2	Ethane, diazo-	56	C2H4N2	001117-96-0	9
3	2-Butene, (E)-	56	C4H8	000624-64-6	9
4	2-Butene	56	C4H8	000107-01-7	9



Library Search Compound Report

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Acq On : 30 Nov 2001 10:13 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
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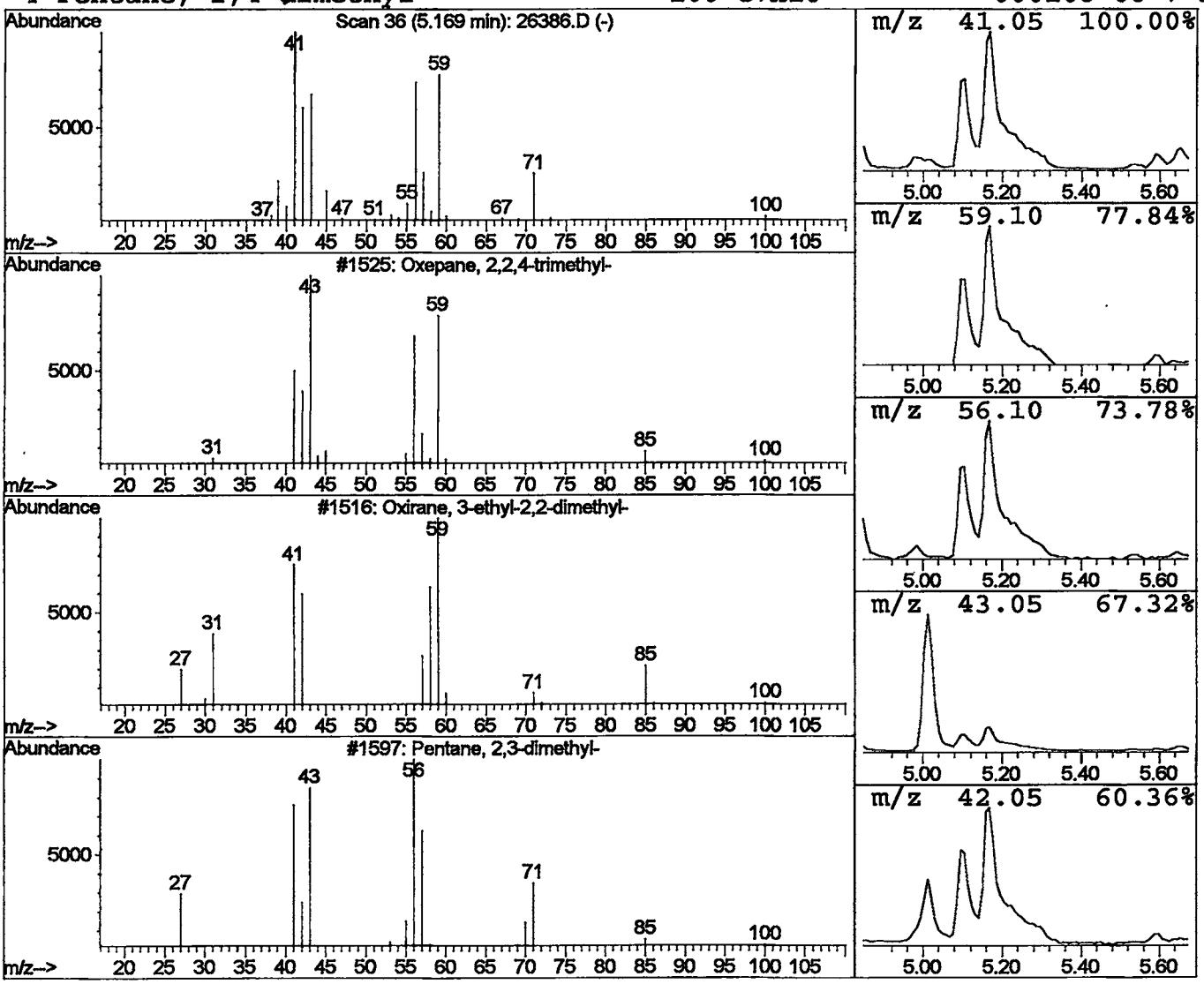
****

Peak Number 3 Oxepane, 2,2,4-trimethyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.17	2.13 ng/uL	949598	1,4-Dichlorobenzene-d4	11.89

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Oxepane, 2,2,4-trimethyl-	100	C6H12O	023120-44-7	47
2	Oxirane, 3-ethyl-2,2-dimethyl-	100	C6H12O	001192-22-9	38
3	Pentane, 2,3-dimethyl-	100	C7H16	000565-59-3	35
4	Pentane, 2,4-dimethyl-	100	C7H16	000108-08-7	32



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Misc :
MS Integration Params: LSCINT.P

Vial: 12
Operator: GALOS
Inst : GC/MS Ins
Multiplr: 1.00

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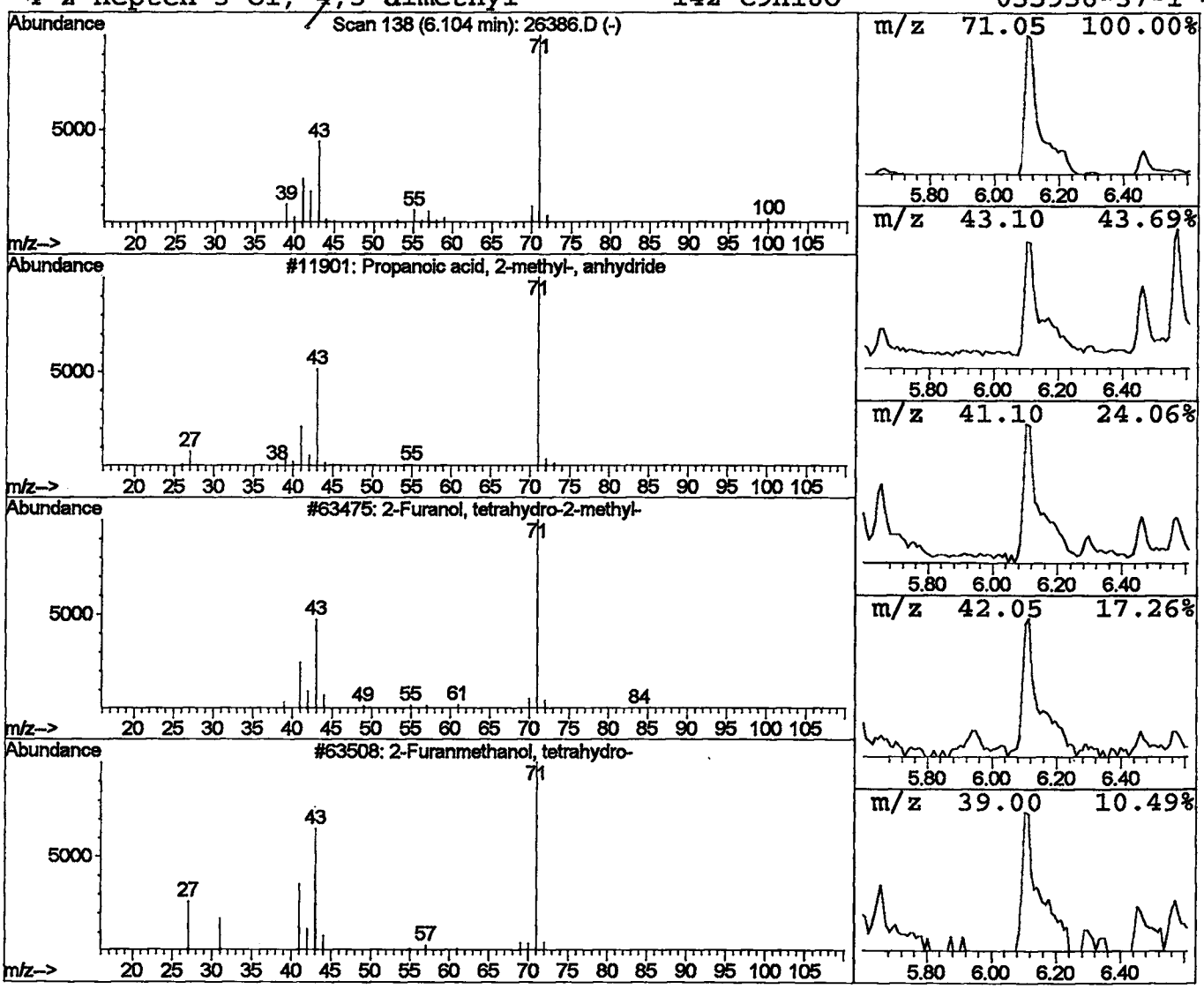
Peak Number 4 Propanoic acid, 2-methyl-, anh Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.10	1.08 ng/uL	480759	1,4-Dichlorobenzene-d4	11.89

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Propanoic acid, 2-methyl-, anhydrid	158	C8H14O3	000097-72-3	59
2	2-Furanol, tetrahydro-2-methyl-	102	C5H10O2	007326-46-7	43
3	2-Furanmethanol, tetrahydro-	102	C5H10O2	000097-99-4	40
4	2-Hepten-3-ol, 4,5-dimethyl-	142	C9H18O	055956-37-1	40



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Multiplr: 1.00

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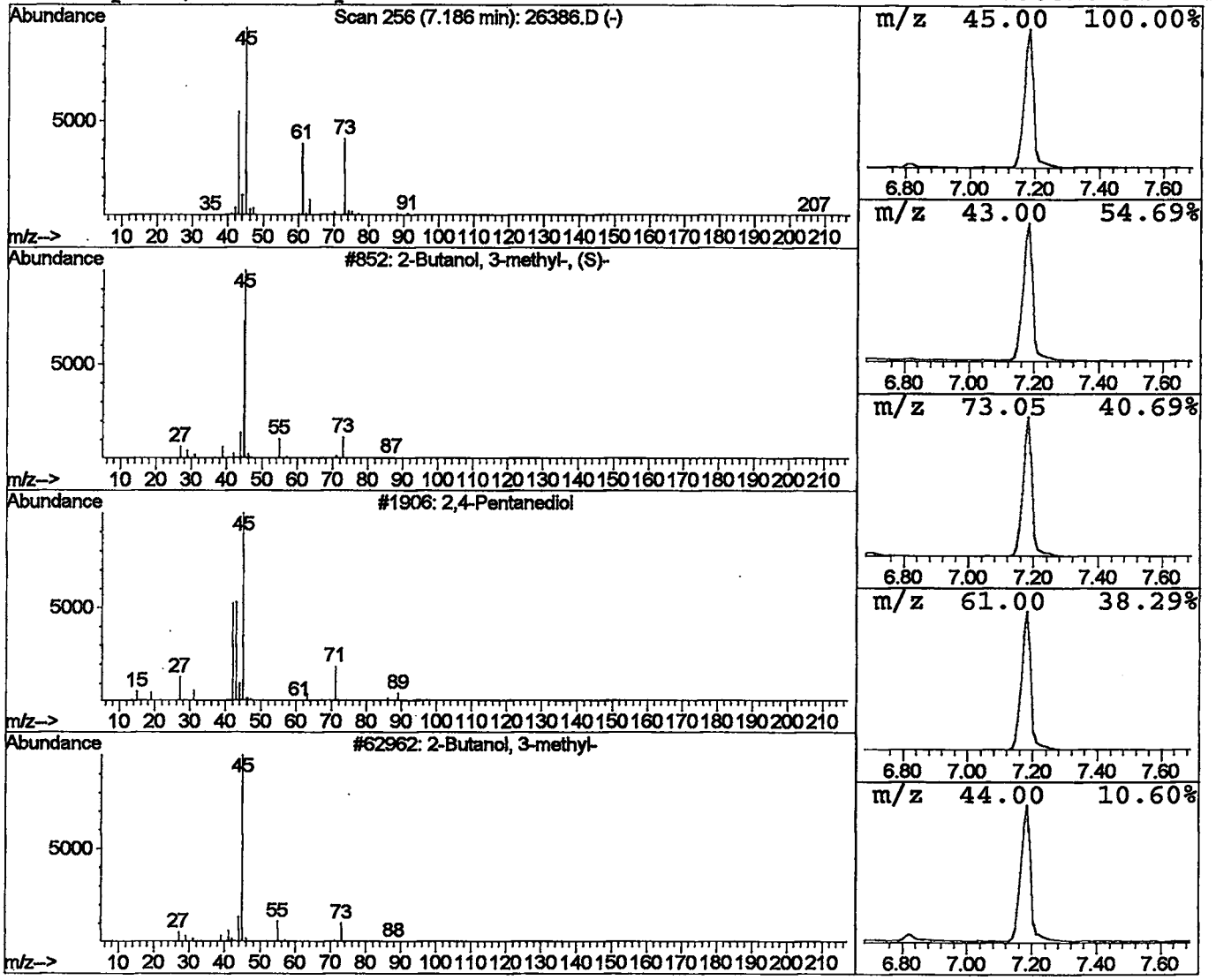
Library : C:\DATABASE\NBS75K.L

****

Peak Number 5 2-Butanol, 3-methyl-, (S)- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.19	10.04 ng/uL	4480040	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	2,4-Pentanediol	104	C5H12O2	000625-69-4	33
3	2-Butanol, 3-methyl-	88	C5H12O	000598-75-4	9
4	Propane, 2-ethoxy-	88	C5H12O	000625-54-7	9



Library Search Compound Report

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

Vial: 12
Operator: GALOS
Inst : GC/MS Ins
Multiplr: 1.00

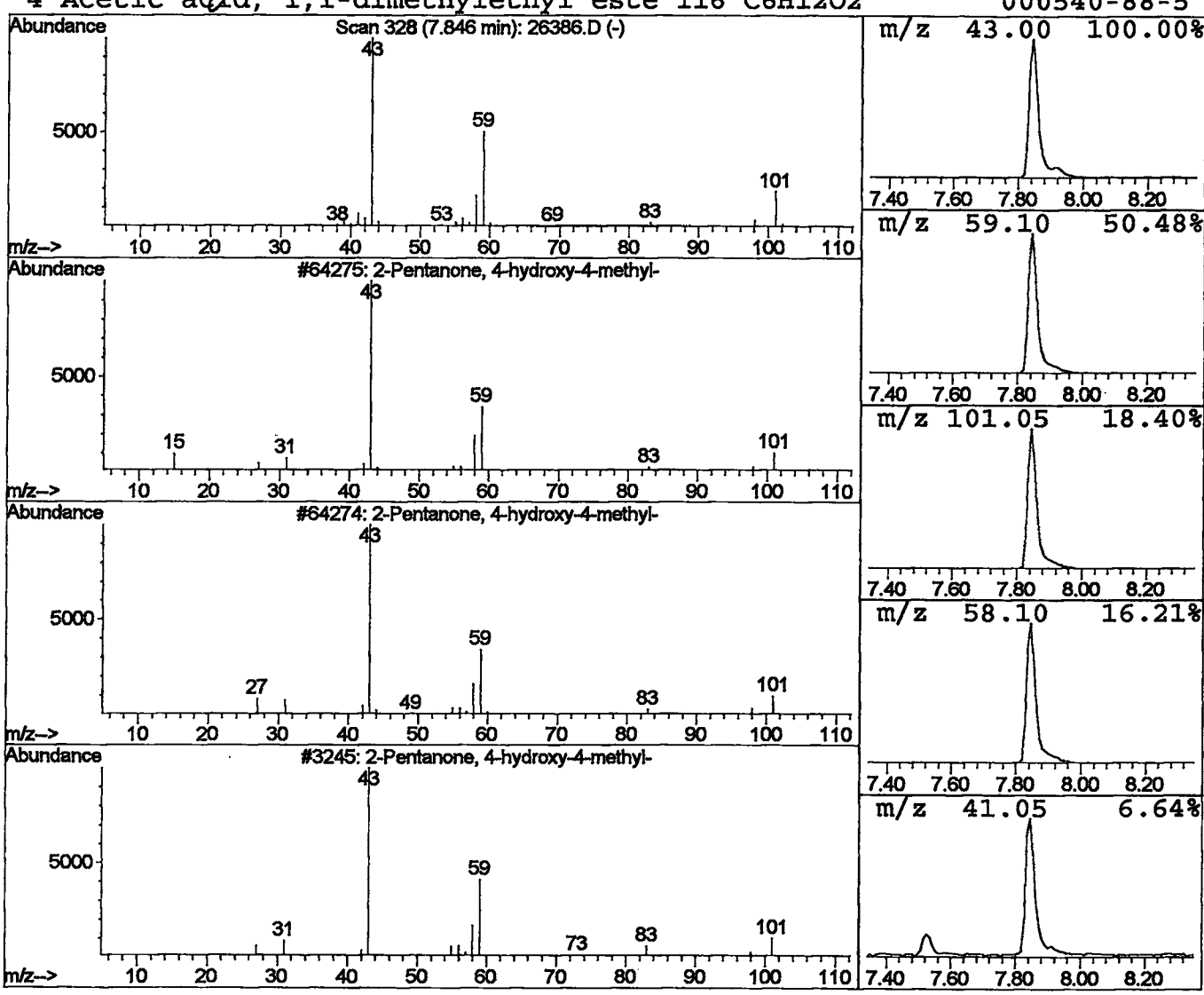
Quant Method : C:\HPCHEM\1\METHODS\NP112701.M (RTE Integrator)
Title : 208 625/TCLP calibration table
Library : C:\DATABASE\NBS75K.L

****

Peak Number 6 2-Pentanone, 4-hydroxy-4-methy Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.85	5.17 ng/uL	2304210	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	53
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	43
4	Acetic acid, 1,1-dimethylethyl este	116	C6H12O2	000540-88-5	28



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MS Integration Params: LSCINT.P

Vial: 12
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Inst : GC/MS Ins
Multiplr: 1.00

Quant Method : C:\HPCHEM\1\METHODS\NP112701.M (RTE Integrator)
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Library : C:\DATABASE\NBS75K.L

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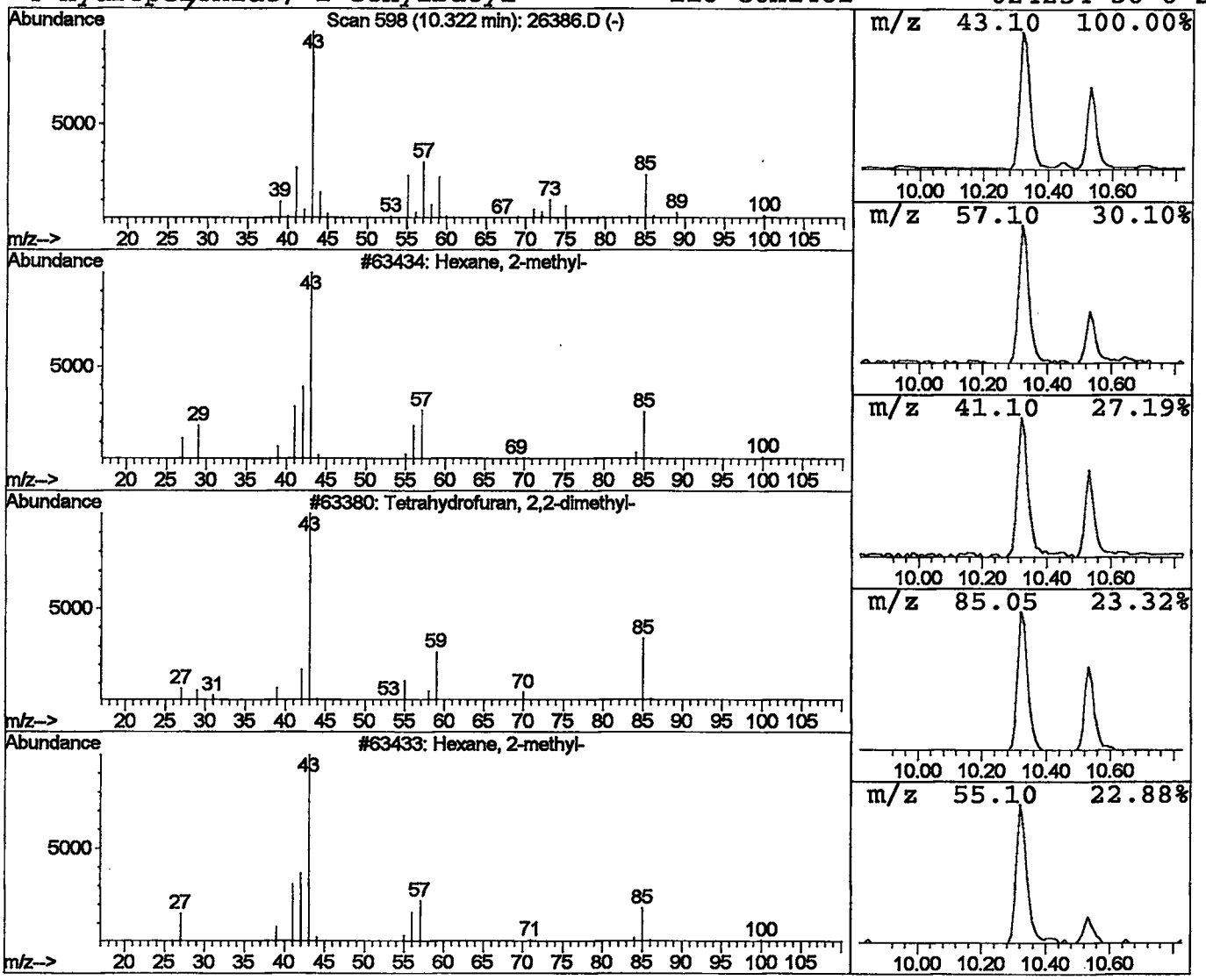
Peak Number 7 Hexane, 2-methyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.32	1.69 ng/uL	753211	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	Tetrahydrofuran, 2,2-dimethyl-	100	C6H12O	001003-17-4	33
3	Hexane, 2-methyl-	100	C7H16	000591-76-4	25
4	Hydroperoxide, 1-ethylbutyl	118	C6H14O2	024254-56-6	23



Library Search Compound Report

Data File : C:\HPCHEM\1\DATA\NOV3001\26386.D
Acq On : 30 Nov 2001 10:13 pm
Sample : gc/ms unknown 11/3
Misc :
MS Integration Params: LSCINT.P

Vial: 12
Operator: GALOS
Inst : GC/MS Ins
Multiplr: 1.00

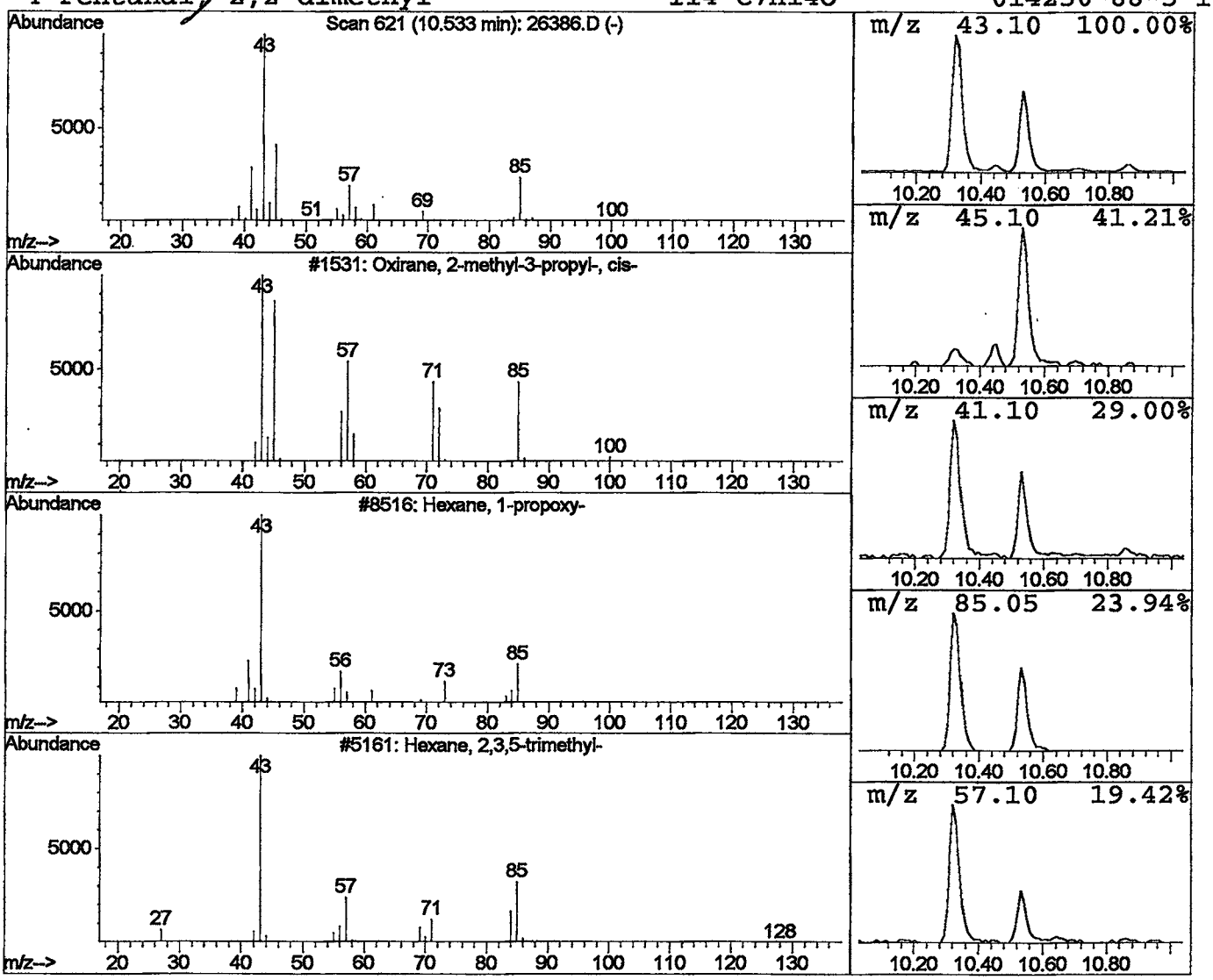
Quant Method : C:\HPCHEM\1\METHODS\NP112701.M (RTE Integrator)
Title : 208 625/TCLP calibration table
Library : C:\DATABASE\NBS75K.L

****

Peak Number 8 Oxirane, 2-methyl-3-propyl-, c Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.53	0.86 ng/uL	384399	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	36
2	Hexane, 1-propoxy-	144	C9H20O	053685-78-2	23
3	Hexane, 2,3,5-trimethyl-	128	C9H20	001069-53-0	16
4	Pentanal, 2,2-dimethyl-	114	C7H14O	014250-88-5	16



Library Search Compound Report

Data File : C:\HPCHEM\1\DATA\NOV3001\26386.D
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Sample : gc/ms unknown 11/3
Misc :
MS Integration Params: LSCINT.P

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Inst : GC/MS Ins
Multiplr: 1.00

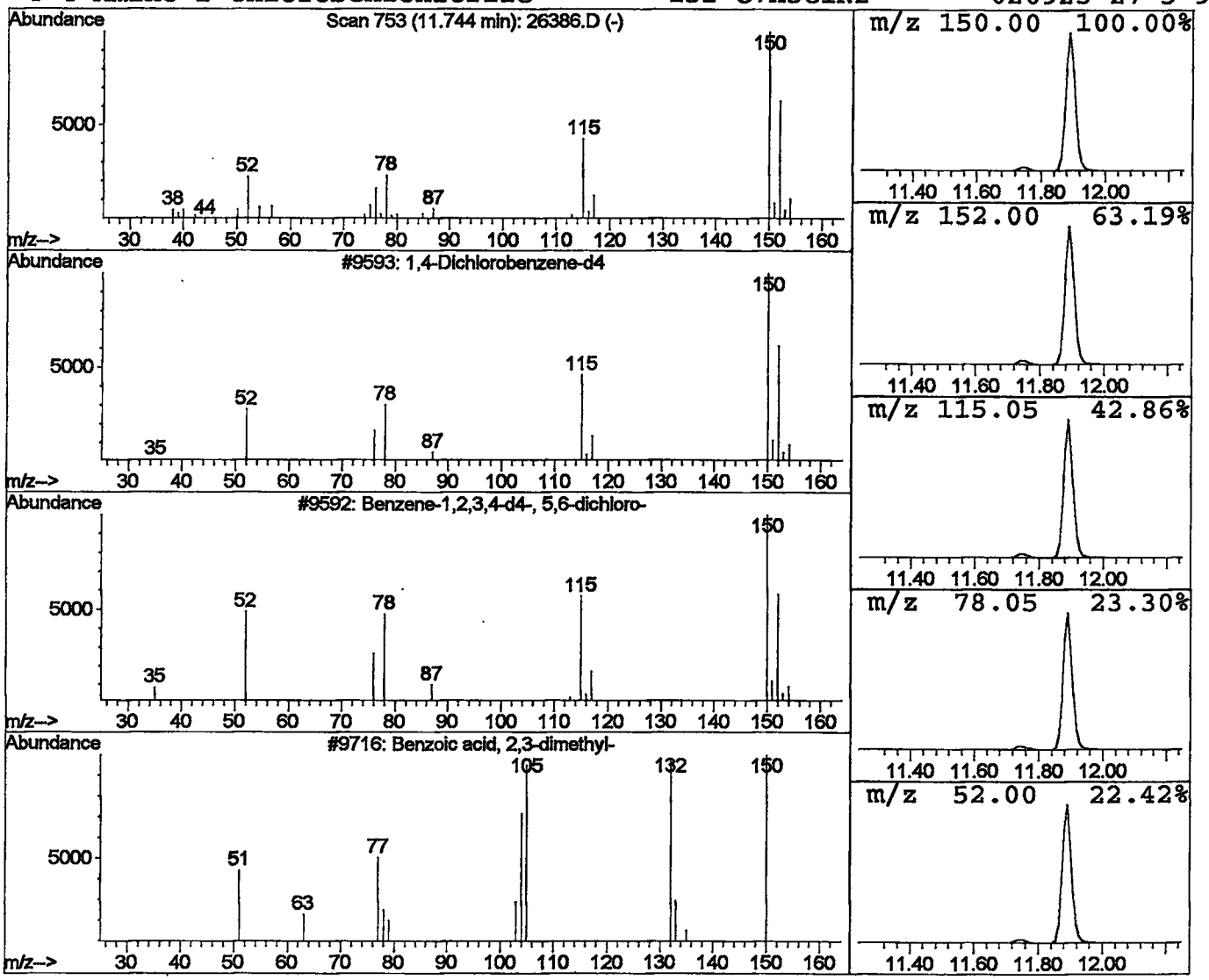
Quant Method : C:\HPCHEM\1\METHODS\NP112701.M (RTE Integrator)
Title : 208 625/TCLP calibration table
Library : C:\DATABASE\NBS75K.L

****

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.74	0.55 ng/uL	244101	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	96
2	Benzene-1,2,3,4-d4-, 5,6-dichloro-	150	C6Cl2D4	002199-69-1	53
3	Benzoic acid, 2,3-dimethyl-	150	C9H10O2	000603-79-2	9
4	4-Amino-2-chlorobenzonitrile	152	C7H5ClN2	020925-27-3	9



Tentatively Identified Compound (LSC) summary

Operator ID: GALOS Date Acquired: 30 Nov 2001 10:13 pm
Data File: C:\HPCHEM\1\DATA\NOV3001\26386.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
Hydrogen azide	5.01	1.3	ng/uL	580103	ISTD01	11.89	8921270	20.0
Cyclopropane, (1-met	5.10	0.8	ng/uL	365882	ISTD01	11.89	8921270	20.0
Oxepane, 2,2,4-trime	5.17	2.1	ng/uL	949598	ISTD01	11.89	8921270	20.0
Propanoic acid, 2-me	6.10	1.1	ng/uL	480759	ISTD01	11.89	8921270	20.0
2-Butanol, 3-methyl-	7.19	10.0	ng/uL	4480040	ISTD01	11.89	8921270	20.0
2-Pentanone, 4-hydro	7.85	5.2	ng/uL	2304210	ISTD01	11.89	8921270	20.0
Hexane, 2-methyl-	10.32	1.7	ng/uL	753211	ISTD01	11.89	8921270	20.0
Oxirane, 2-methyl-3-	10.53	0.9	ng/uL	384399	ISTD01	11.89	8921270	20.0
1,4-Dichlorobenzene-	11.74	0.5	ng/uL	244101	ISTD01	11.89	8921270	20.0

26386.D NP112701.M Tue Dec 04 10:38:38 2001