User: Galos, Marie
Westchester Dept. of Labs & Research
Date: 12/04/2001 Time: 15:13:52

Sample: AD26389
Analysis: $GCMS GCMS Scan For Unknowns
Units: ug
Started: 12/4/01 15:13 Ended: 12/4/01 15:13 Analyst: MG
Page: 1

	Component Name	Viol	Result
1	Other unknown compounds		see comment

Comments for Sample AD26389 - Analysis $GCMS

Westchester Dept. of Labs & Research
User: Galos, Marie
Date: 12-04-2001 Time: 15:14:03

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\26389.D Vial: 6
Acq On : 30 Nov 2001 4:46 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 14:53 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	1410369	20.00	ng/uL	-0.04
19) Naphthalene-d8	15.50	136	5451984	20.00	ng/uL	-0.04
31) Acenaphthene-d10	20.78	164	2473650	20.00	ng/uL	-0.04
49) Phenanthrene-d10	25.19	188	3520255	20.00	ng/uL	-0.05
59) Chrysene-d12	33.19	240	2203164	20.00	ng/uL	-0.05
68) Perylene-d12	37.27	264	1563013	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	11.19	99	1035	0.01	ng/uL	0.07
Spiked Amount 75.000	Range 19 - 32		Recovery =	0.01%#		
6) 2-Chlorophenol-d4	11.88	132	1099	0.01	ng/uL	0.46
Spiked Amount 75.000	Range 34 - 84		Recovery =	0.01%#		
7) 1,2-Dichlorobenzene-d4	12.40	152	14074	0.24	ng/uL	-0.05
Spiked Amount 50.000	Range 26 - 73		Recovery =	0.48%#		
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	4770	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	287	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

Quantitation Report (QT Reviewed)

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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	769	N.D.	
56) Anthracene	25.19	178	769	N.D.	
57) Di-n-butylphthalate	27.18	149	3305	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	33.18	228	4820	N.D.	
66) Bis(2-ethylhexyl)phthalate	33.45	149	6220	N.D.	
67) Chrysene	33.18	228	4820	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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	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	5353	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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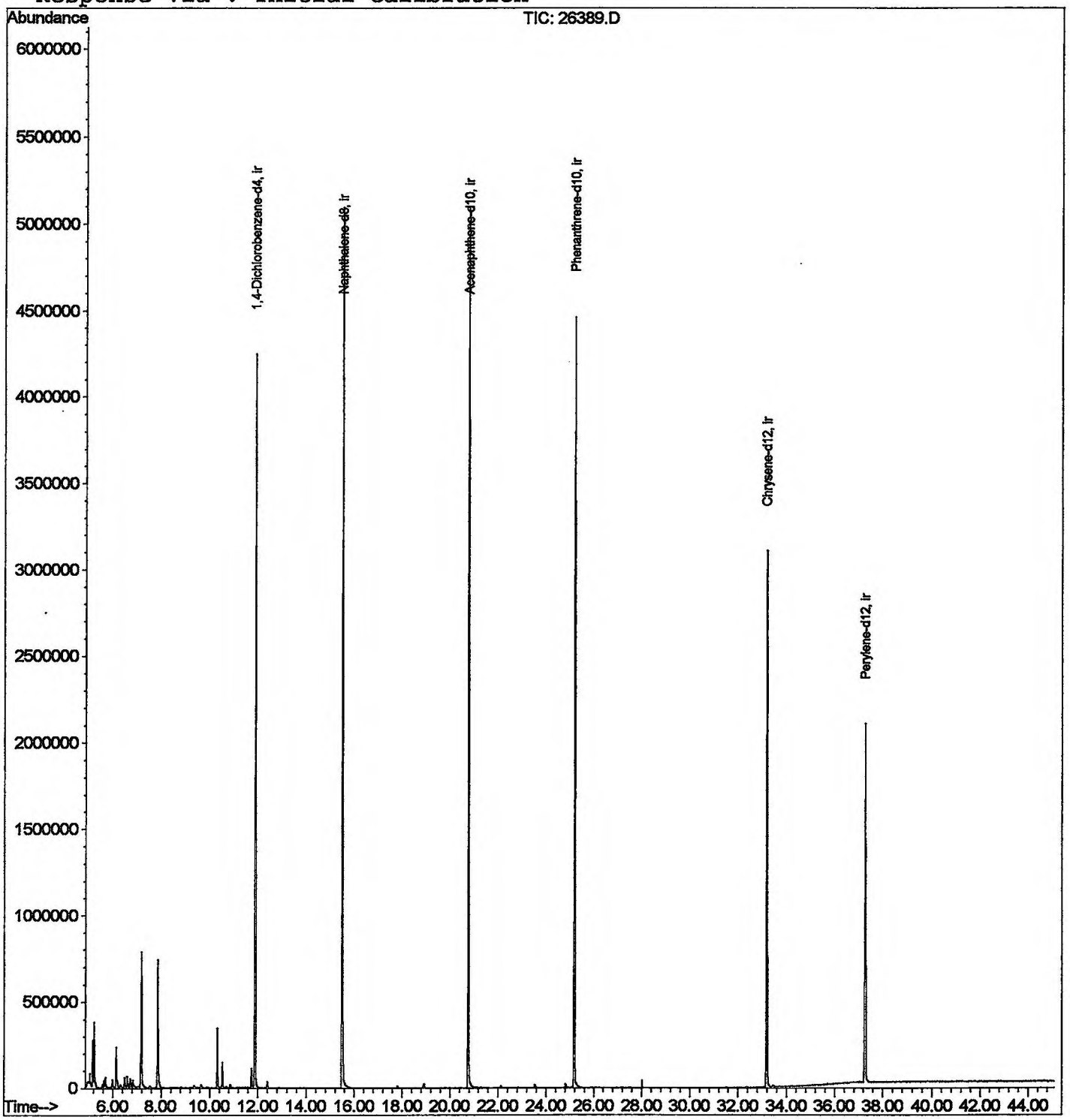


LSC Area Percent Report

Data File : C:\HPCHEM\1\DATA\NOV3001\26389.D Vial: 6
Acq On : 30 Nov 2001 4:46 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

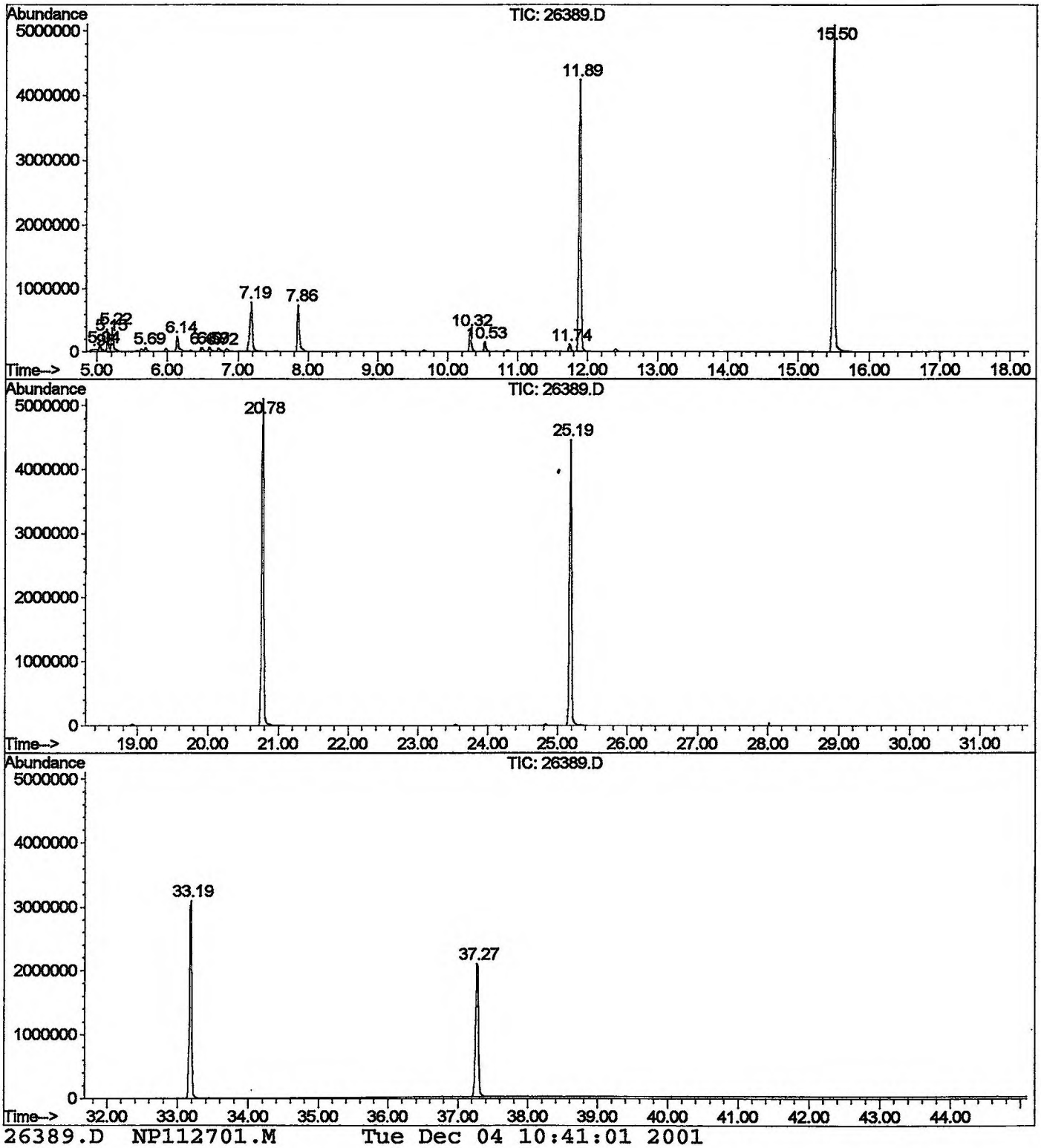
peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	peak area	peak % max.	% of total
1	4.977	5	15	19	rBV	30994	167333	1.53%	0.282%
2	5.041	19	22	31	rVV2	77354	208839	1.91%	0.351%
3	5.151	31	34	38	rVV	273132	445903	4.08%	0.750%
4	5.216	38	41	54	rVB	380409	741706	6.79%	1.248%
5	5.692	89	93	102	rVB	61551	130502	1.19%	0.220%
6	6.142	138	142	155	rVB	236370	502028	4.60%	0.845%
7	6.490	175	180	187	rBV	57784	123265	1.13%	0.207%
8	6.591	187	191	201	rVB	65229	140686	1.29%	0.237%
9	6.719	201	205	213	rBV2	50085	119134	1.09%	0.200%
10	7.187	248	256	266	rBV	785232	1887330	17.28%	3.175%
11	7.857	325	329	342	rBV	741237	1497545	13.71%	2.519%
12	10.323	593	598	608	rBV	347601	656622	6.01%	1.105%
13	10.534	616	621	629	rBV	150237	285903	2.62%	0.481%
14	11.745	749	753	761	rBV	113503	229246	2.10%	0.386%
15	11.891	763	769	781	rVB	4243968	8505172	77.86%	14.309%
16	15.504	1156	1163	1180	rBV	5094031	10923701	100.00%	18.377%
17	20.777	1731	1738	1758	rBV	5102113	10635949	97.37%	17.893%
18	25.188	2212	2219	2230	rBV2	4463894	9541093	87.34%	16.051%
19	33.194	3084	3092	3104	rBV2	3103934	7035184	64.40%	11.836%
20	37.274	3529	3537	3552	rBV2	2077700	5663516	51.85%	9.528%

Sum of corrected areas: 59440657

26389.D NP112701.M Tue Dec 04 10:41:00 2001

LSC Report - Integrated Chromatogram

File : C:\HPCHEM\1\DATA\NOV3001\26389.D
Operator : GALOS
Acquired : 30 Nov 2001 4:46 pm using AcqMethod NP112701
Instrument : GC/MS Ins
Sample Name: gc/ms unknown 11/3
Misc Info :
Vial Number: 6
Quant File :NP112701.RES (RTE Integrator)



26389.D NP112701.M Tue Dec 04 10:41:01 2001

Library Search Compound Report

Data File : C:\HPCHEM\1\DATA\NOV3001\26389.D
Acq On : 30 Nov 2001 4:46 pm
Sample : gc/ms unknown 11/3
Misc :
MS Integration Params: LSCINT.P

Vial: 6
Operator: GALOS
Inst : GC/MS Ins
Multiplr: 1.00

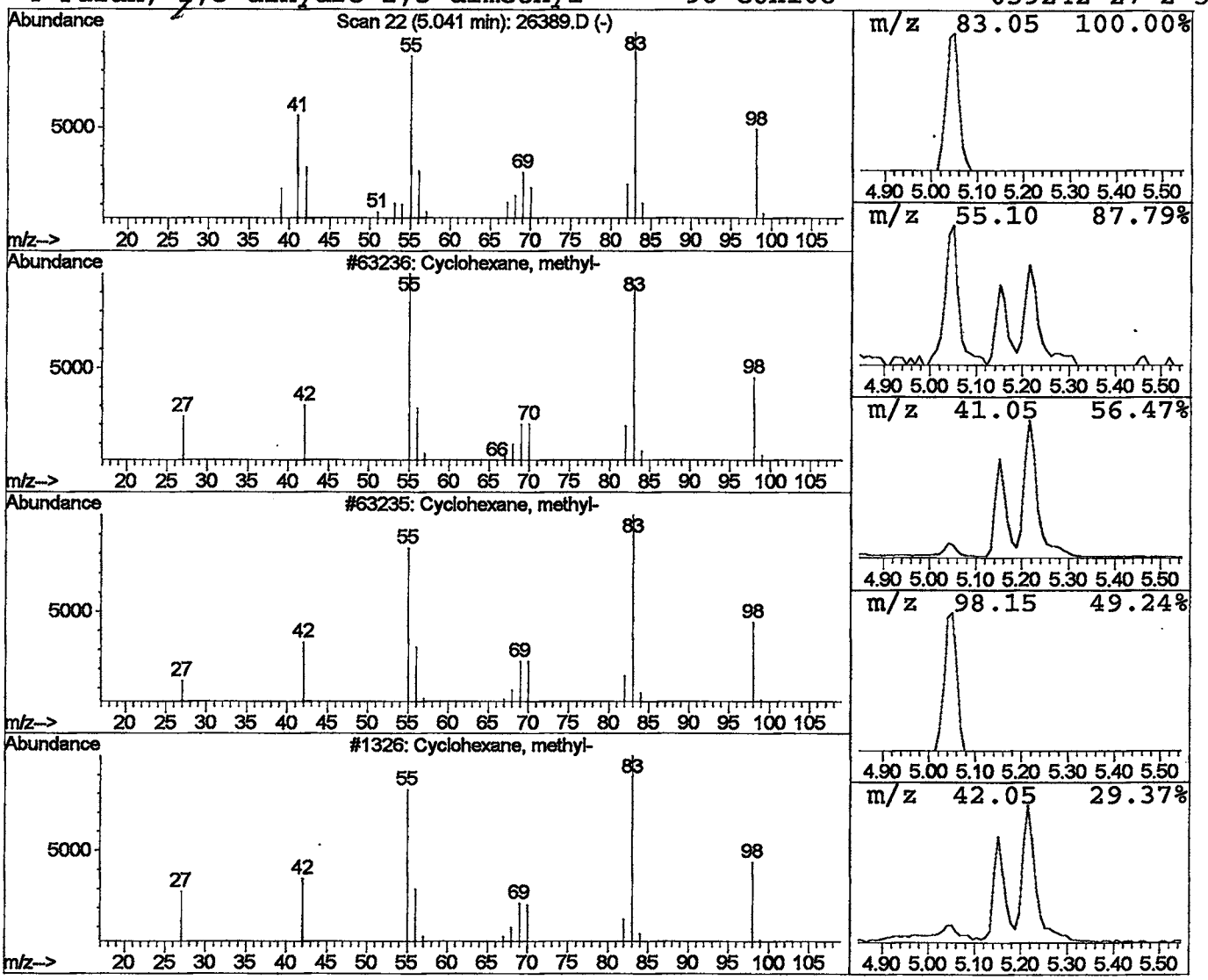
Quant Method : C:\HPCHEM\1\METHODS\NP112701.M (RTE Integrator)
Title : 208 625/TCLP calibration table
Library : C:\DATABASE\NBS75K.L

****

Peak Number 1 Cyclohexane, methyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.04	0.49 ng/uL	208839	1,4-Dichlorobenzene-d4	11.89

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclohexane, methyl-	98	C7H14	000108-87-2	97
2	Cyclohexane, methyl-	98	C7H14	000108-87-2	94
3	Cyclohexane, methyl-	98	C7H14	000108-87-2	91
4	Furan, 2,5-dihydro-2,5-dimethyl-	98	C6H10O	059242-27-2	59



Library Search Compound Report

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Acq On : 30 Nov 2001 4:46 pm
Sample : gc/ms unknown 11/3
Misc :
MS Integration Params: LSCINT.P

Vial: 6
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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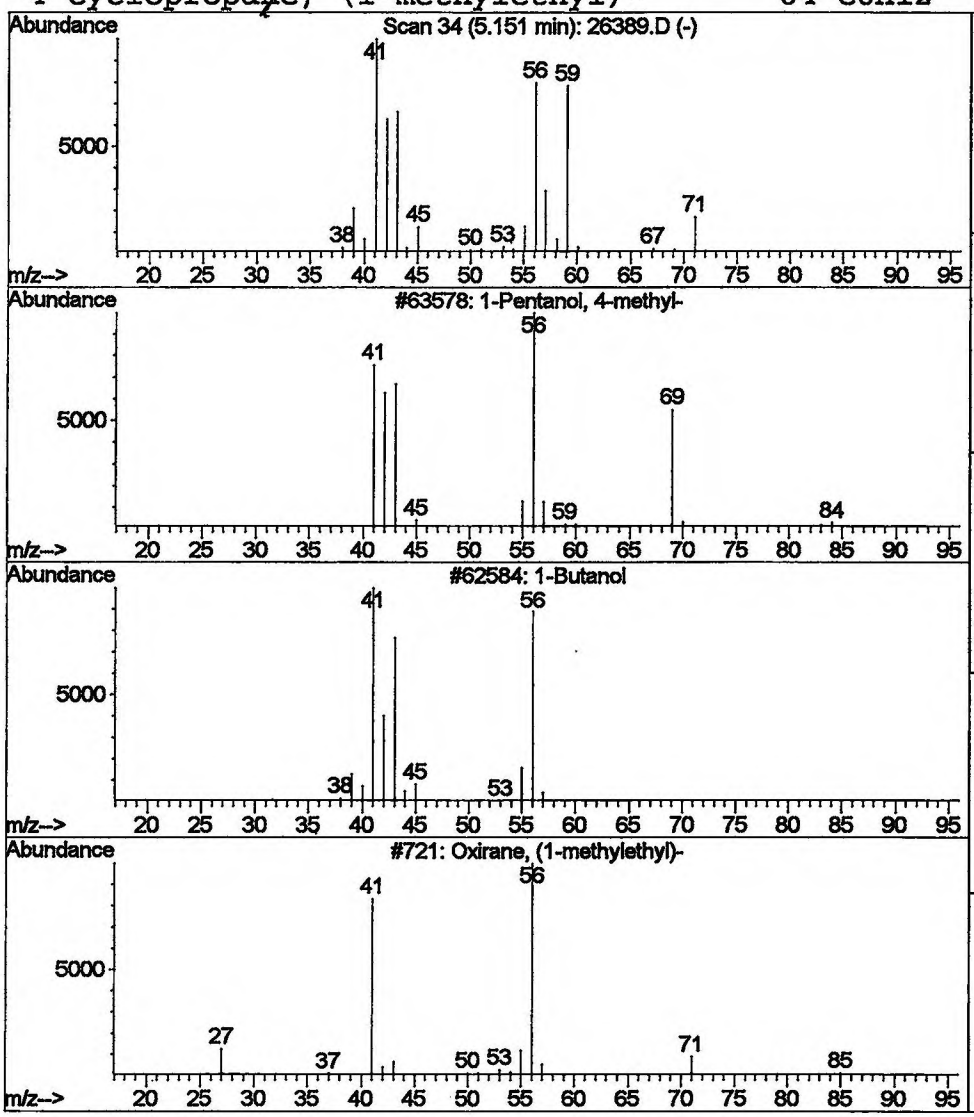
****

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.15	1.05 ng/uL	445903	1,4-Dichlorobenzene-d4	11.89

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Pentanol, 4-methyl-	102	C6H14O	000626-89-1	25
2	1-Butanol	74	C4H10O	000071-36-3	23
3	Oxirane, (1-methylethyl)-	86	C5H10O	001438-14-8	12
4	Cyclopropane, (1-methylethyl)-	84	C6H12	003638-35-5	10



Library Search Compound Report

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Acq On : 30 Nov 2001 4:46 pm
Sample : gc/ms unknown 11/3
Misc :
MS Integration Params: LSCINT.P

Vial: 6
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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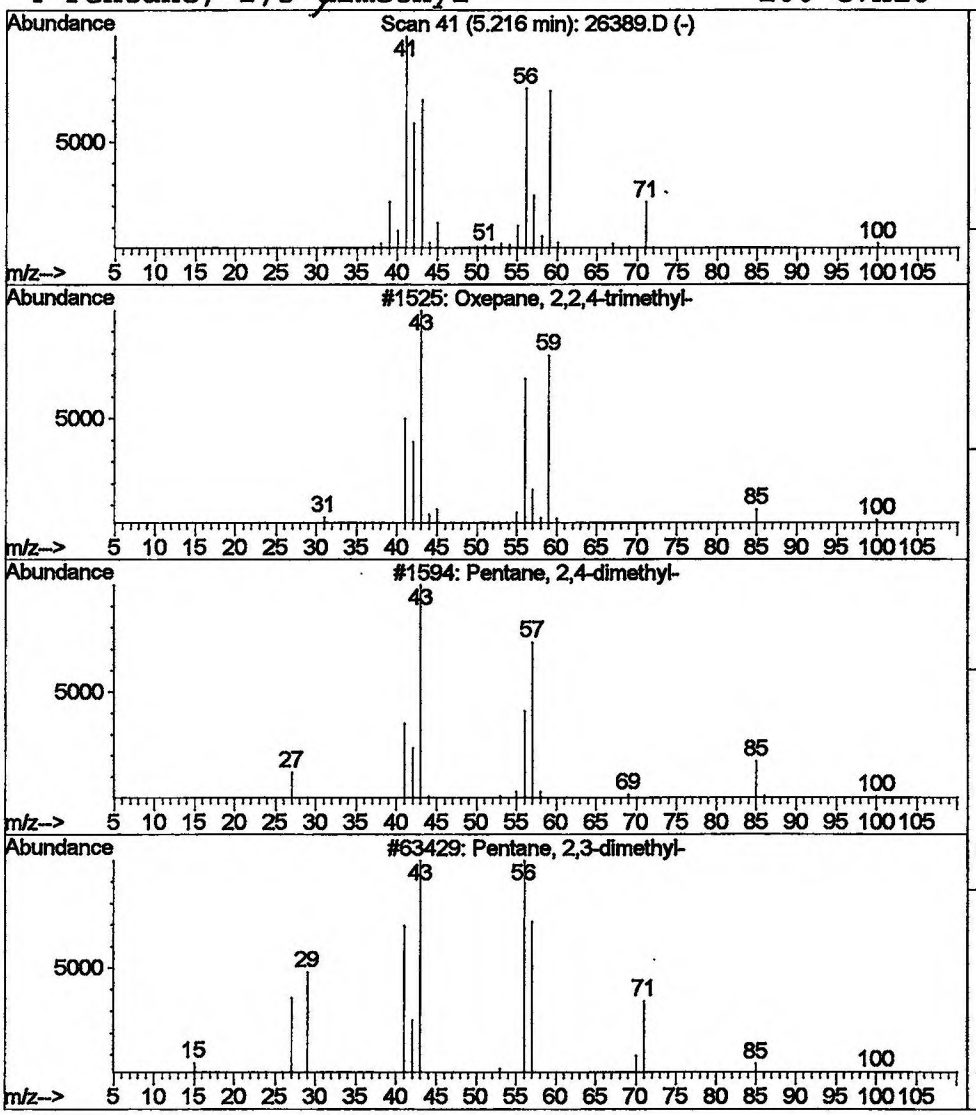
****

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.22	1.74 ng/uL	741706	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	64
2	Pentane, 2,4-dimethyl-	100	C7H16	000108-08-7	38
3	Pentane, 2,3-dimethyl-	100	C7H16	000565-59-3	35
4	Pentane, 2,3-dimethyl-	100	C7H16	000565-59-3	35



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Acq On : 30 Nov 2001 4:46 pm
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Misc :
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Vial: 6
Operator: GALOS
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Multiplr: 1.00

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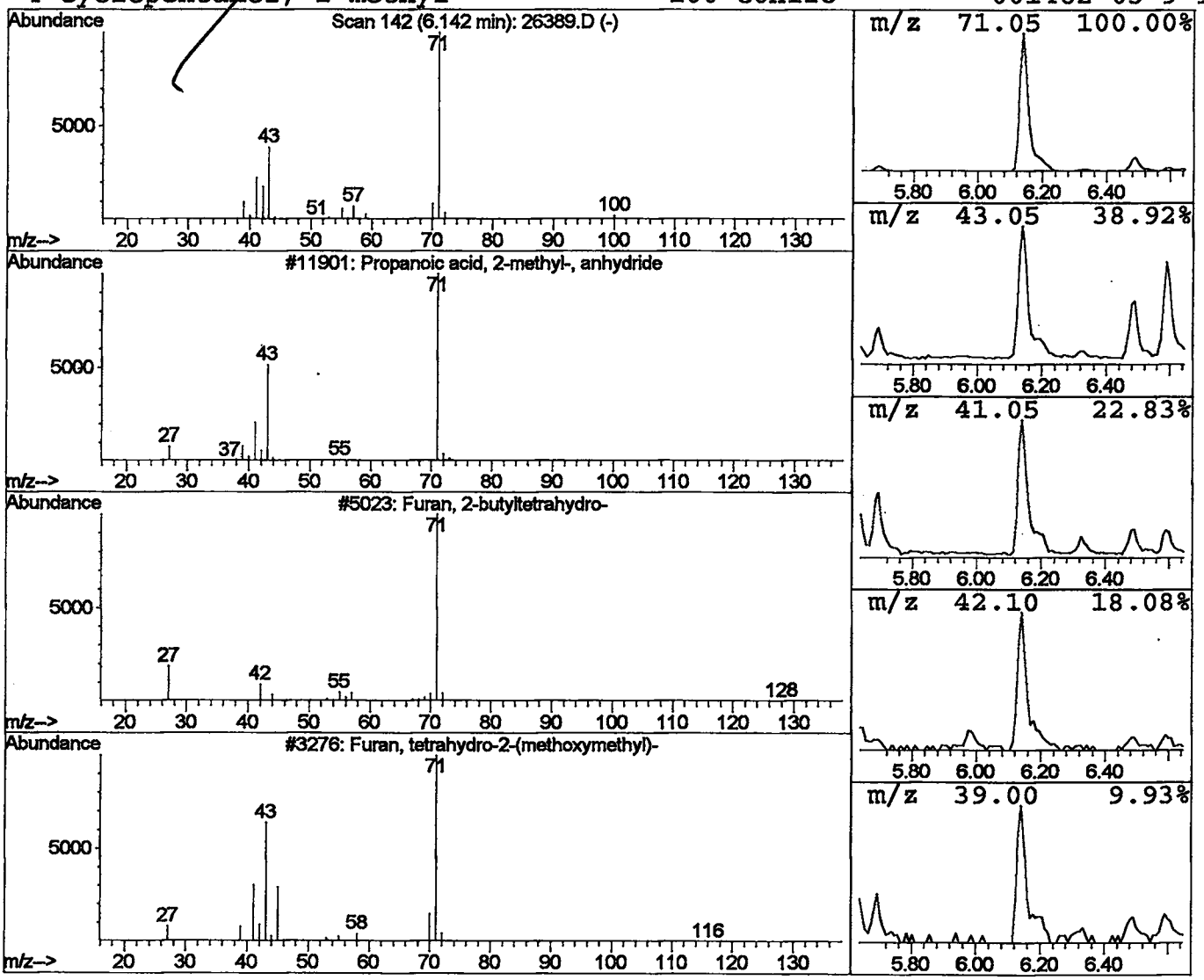
****

Peak Number 4 Propanoic acid, 2-methyl-, anh Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.14	1.18 ng/uL	502028	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	Furan, 2-butyltetrahydro-	128	C8H16O	001004-29-1	39
3	Furan, tetrahydro-2-(methoxymethyl)	116	C6H12O2	019354-27-9	25
4	Cyclopentanol, 1-methyl-	100	C6H12O	001462-03-9	9



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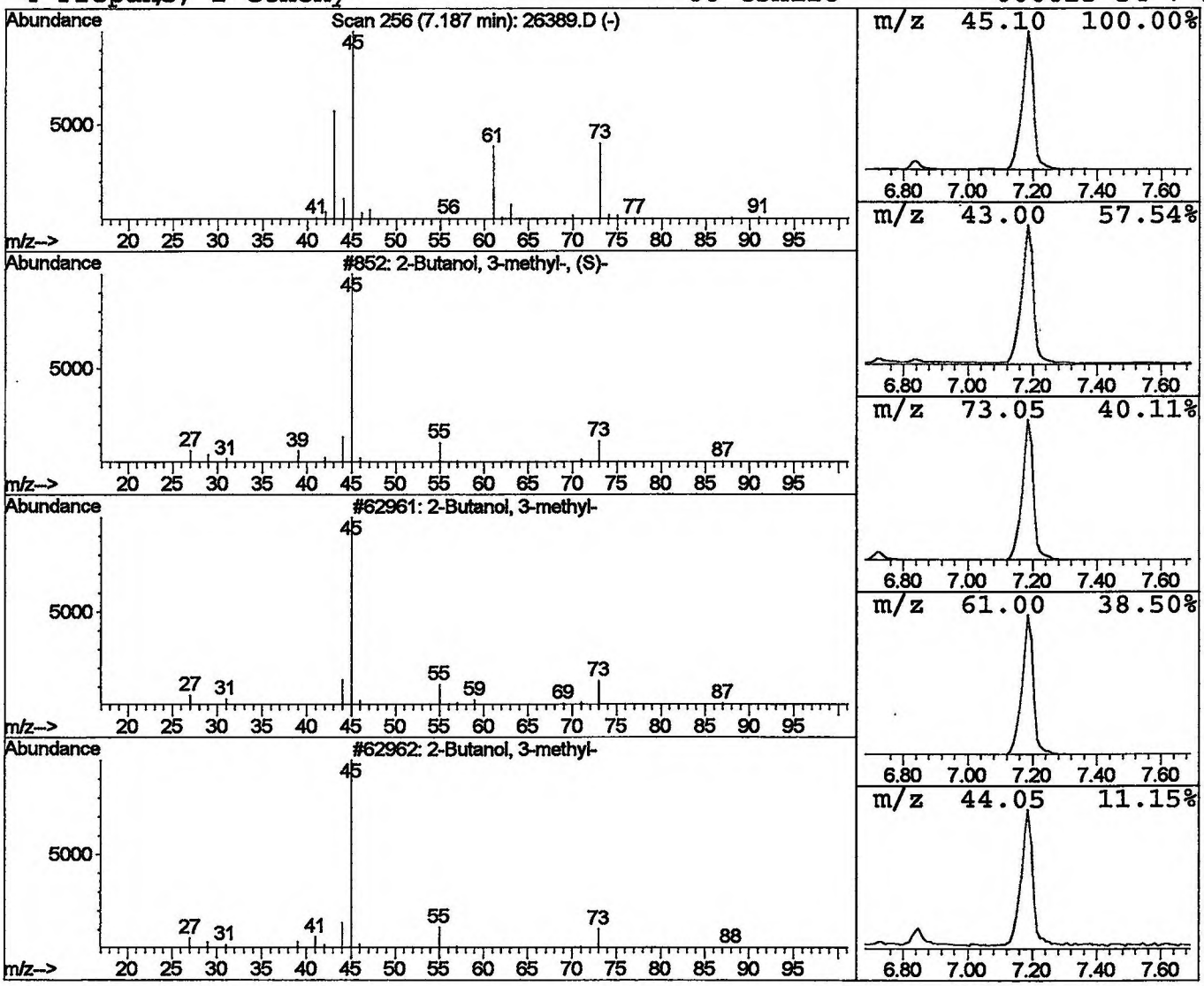
****

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.19	4.44 ng/uL	1887330	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	40
2	2-Butanol, 3-methyl-	88	C5H12O	000598-75-4	9
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: 6
Operator: GALOS
Inst : GC/MS Ins
Multiplr: 1.00

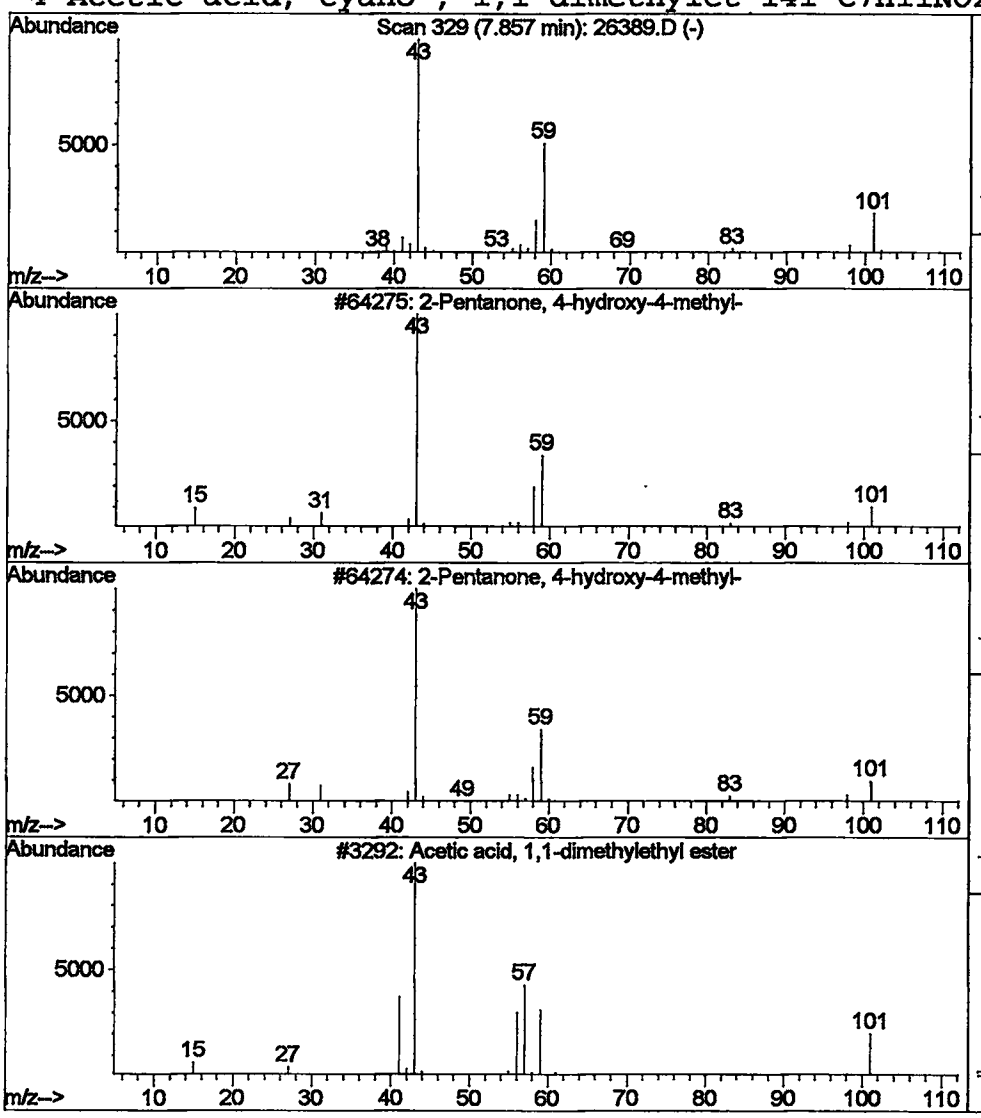
Quant Method : C:\HPCHEM\1\METHODS\NP112701.M (RTE Integrator)
Title : 208 625/TCLP calibration table
Library : C:\DATABASE\NBS75K.L

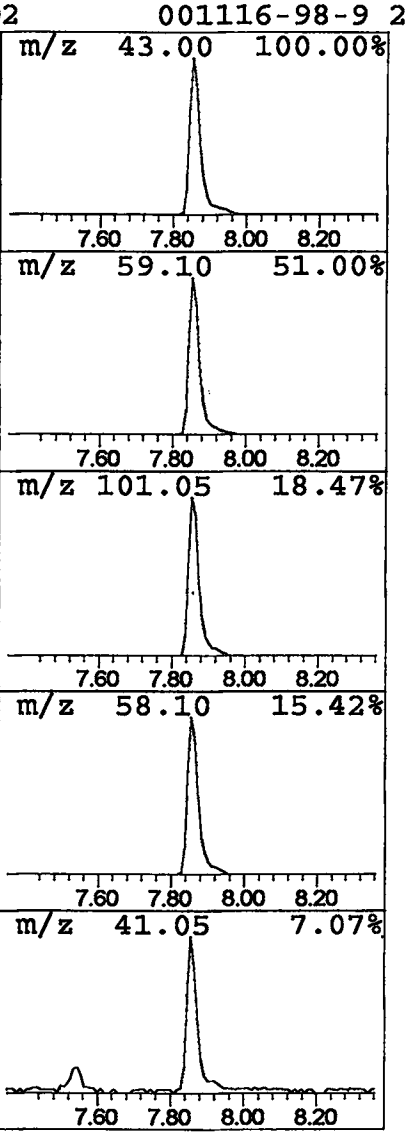
****

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.86	3.52 ng/uL	1497550	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	36
3	Acetic acid, 1,1-dimethylethyl este	116	C6H12O2	000540-88-5	28
4	Acetic acid, cyano-, 1,1-dimethylet	141	C7H11NO2	001116-98-9	23





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

Vial: 6
Operator: GALOS
Inst : GC/MS Ins
Multiplr: 1.00

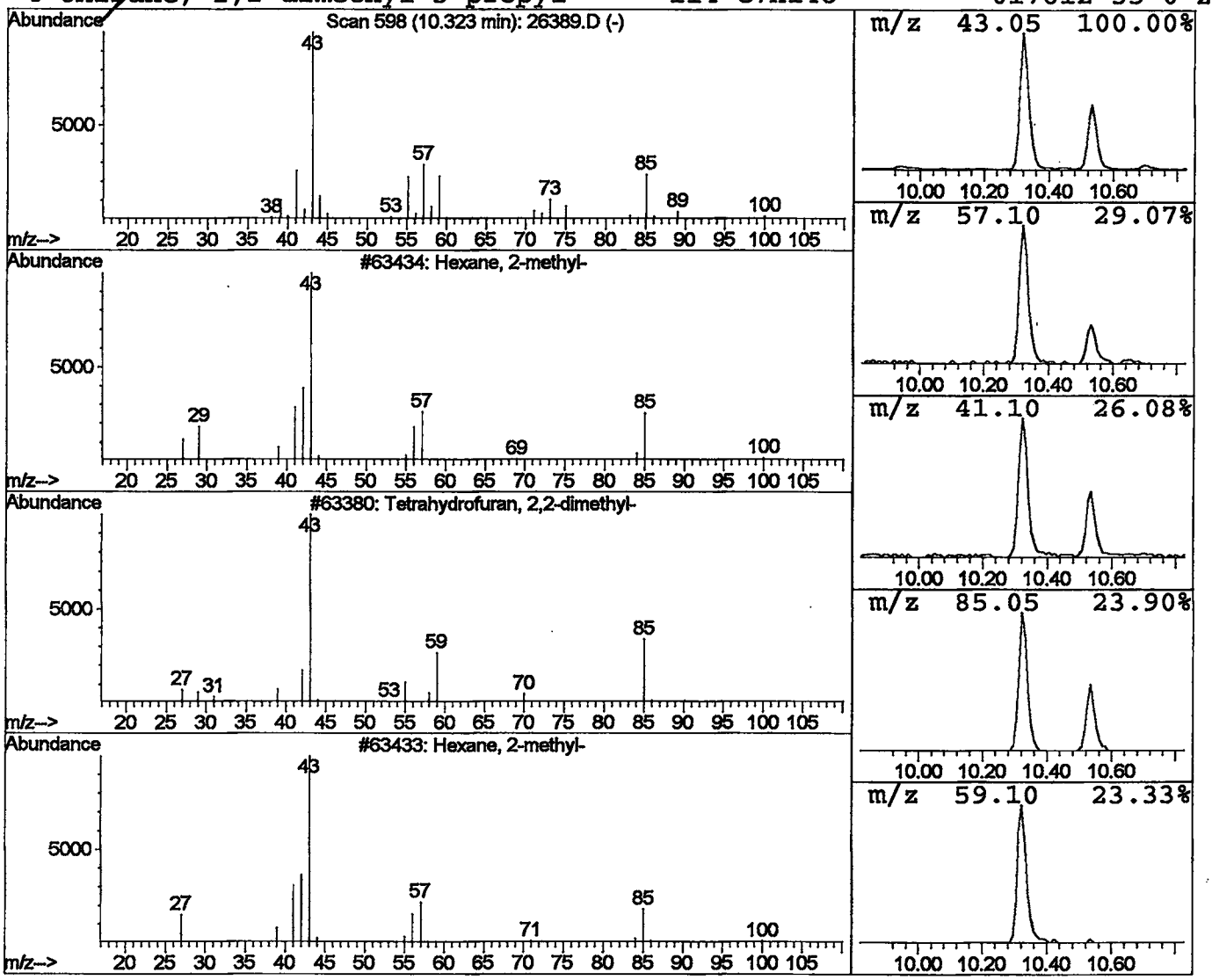
Quant Method : C:\HPCHEM\1\METHODS\NP112701.M (RTE Integrator)
Title : 208 625/TCLP calibration table
Library : C:\DATABASE\NBS75K.L

****

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.32	1.54 ng/uL	656622	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	Oxirane, 2,2-dimethyl-3-propyl-	114	C7H14O	017612-35-0	20



Library Search Compound Report

Data File : C:\HPCHEM\1\DATA\NOV3001\26389.D
Acq On : 30 Nov 2001 4:46 pm
Sample : gc/ms unknown 11/3
Misc :
MS Integration Params: LSCINT.P

Vial: 6
Operator: GALOS
Inst : GC/MS Ins
Multiplr: 1.00

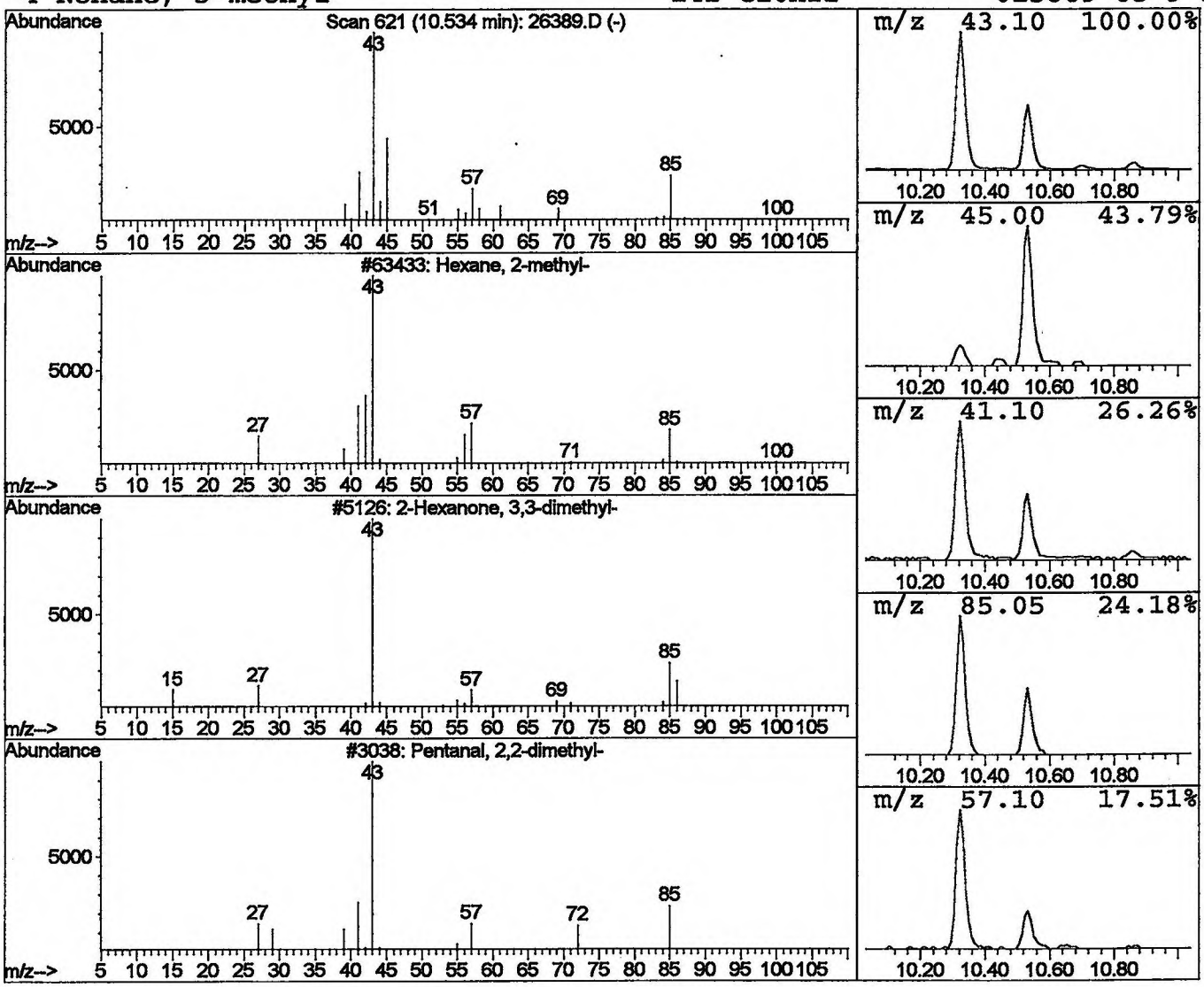
Quant Method : C:\HPCHEM\1\METHODS\NP112701.M (RTE Integrator)
Title : 208 625/TCLP calibration table
Library : C:\DATABASE\NBS75K.L

****

Peak Number 8 Hexane, 2-methyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.53	0.67 ng/uL	285903	1,4-Dichlorobenzene-d4	11.89

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Hexane, 2-methyl-	100	C7H16	000591-76-4	27
2	2-Hexanone, 3,3-dimethyl-	128	C8H16O	026118-38-7	25
3	Pentanal, 2,2-dimethyl-	114	C7H14O	014250-88-5	25
4	Nonane, 5-methyl-	142	C10H22	015869-85-9	17



Library Search Compound Report

Data File : C:\HPCHEM\1\DATA\NOV3001\26389.D
Acq On : 30 Nov 2001 4:46 pm
Sample : gc/ms unknown 11/3
Misc :
MS Integration Params: LSCINT.P

Vial: 6
Operator: GALOS
Inst : GC/MS Ins
Multiplr: 1.00

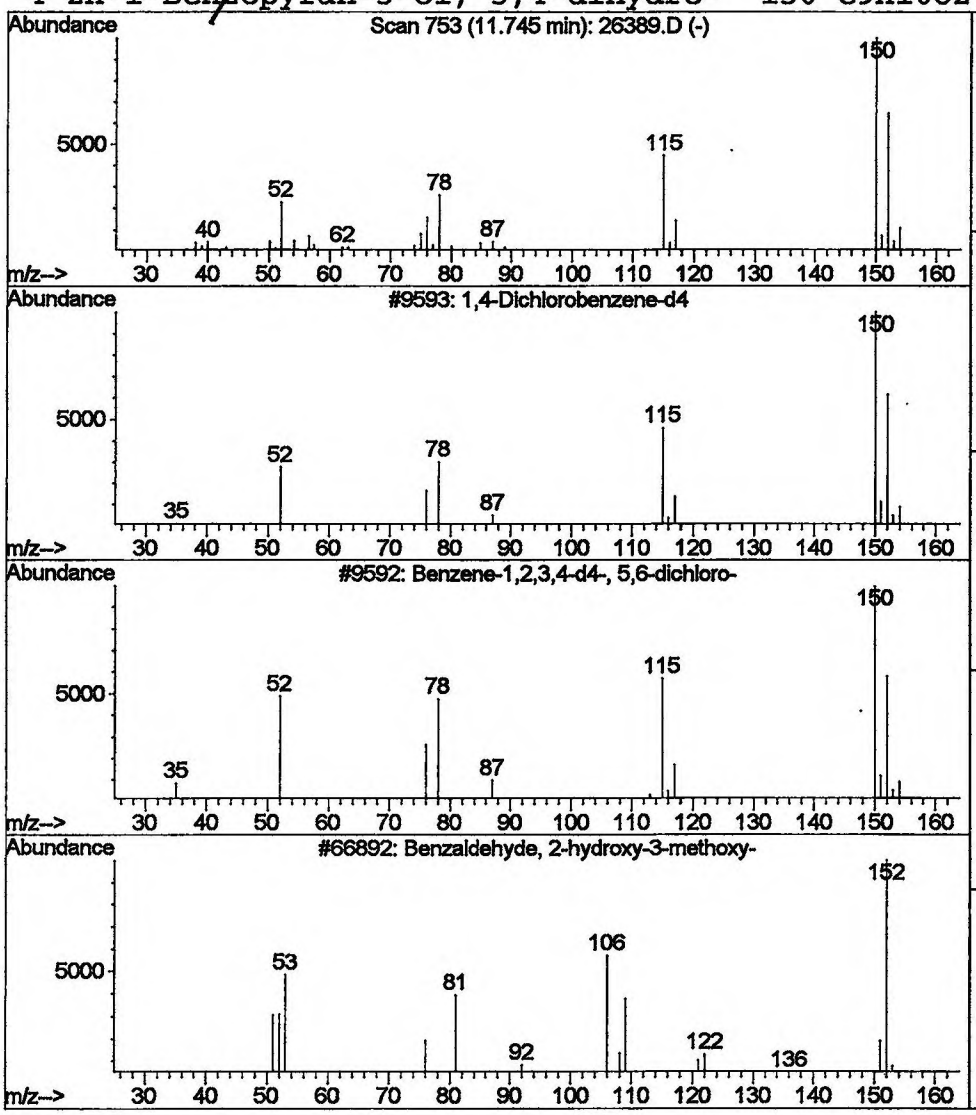
Quant Method : C:\HPCHEM\1\METHODS\NP112701.M (RTE Integrator)
Title : 208 625/TCLP calibration table
Library : C:\DATABASE\NBS75K.L

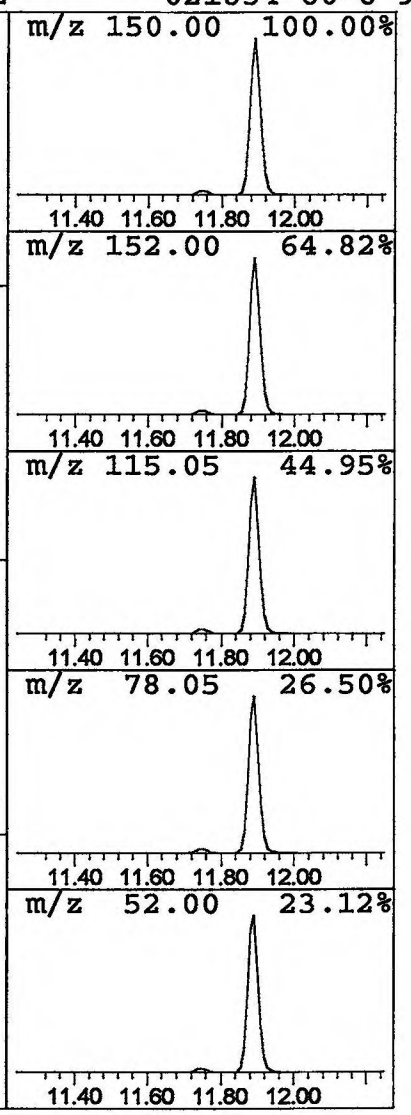
****

Peak Number 9 1,4-Dichlorobenzene-d4 Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.74	0.54 ng/uL	229246	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	94
2	Benzene-1,2,3,4-d4-, 5,6-dichloro-	150	C6Cl2D4	002199-69-1	42
3	Benzaldehyde, 2-hydroxy-3-methoxy-	152	C8H8O3	000148-53-8	9
4	2H-1-Benzopyran-3-ol, 3,4-dihydro-	150	C9H10O2	021834-60-6	9





Tentatively Identified Compound (LSC) summary

Operator ID: GALOS Date Acquired: 30 Nov 2001 4:46 pm
Data File: C:\HPCHEM\1\DATA\NOV3001\26389.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-	5.04	0.5	ng/uL	208839	ISTD01	11.89	8505170	20.0
1-Pentanol, 4-methyl	5.15	1.0	ng/uL	445903	ISTD01	11.89	8505170	20.0
Oxepane, 2,2,4-trime	5.22	1.7	ng/uL	741706	ISTD01	11.89	8505170	20.0
Propanoic acid, 2-me	6.14	1.2	ng/uL	502028	ISTD01	11.89	8505170	20.0
2-Butanol, 3-methyl-	7.19	4.4	ng/uL	1887330	ISTD01	11.89	8505170	20.0
2-Pentanone, 4-hydro	7.86	3.5	ng/uL	1497550	ISTD01	11.89	8505170	20.0
Hexane, 2-methyl-	10.32	1.5	ng/uL	656622	ISTD01	11.89	8505170	20.0
Hexane, 2-methyl-	10.53	0.7	ng/uL	285903	ISTD01	11.89	8505170	20.0
1,4-Dichlorobenzene-	11.74	0.5	ng/uL	229246	ISTD01	11.89	8505170	20.0

26389.D NP112701.M Tue Dec 04 10:41:09 2001