User: Galos, Marie
Westchester Dept. of Labs & Research
Date: 12/10/2001 Time: 14:57:26

Sample: AD26676
Analysis: $GCMS GCMS Scan For Unknowns
Units: ug
Started: 12/10/01 14:57 Ended: 12/10/01 14:57 Analyst: MG
Page: 1

	Component Name	Viol	Result
1	Other unknown compounds		see comment

Comments for Sample AD26676 - Analysis $GCMS

Westchester Dept. of Labs & Research
User: Galos, Marie
Date: 12-10-2001 Time: 14:57:31

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\DEC0601\26676.D Vial: 9
Acq On : 6 Dec 2001 5:28 pm Operator: GALOS
Sample : gc/ms unknown Inst : GC/MS Ins
Misc : Multiplr: 1.00
MS Integration Params: RTEINT.P
Quant Time: Dec 7 10:14 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	999296	20.00	ng/uL	-0.04
19) Naphthalene-d8	15.50	136	3761156	20.00	ng/uL	-0.05
31) Acenaphthene-d10	20.77	164	1716951	20.00	ng/uL	-0.05
49) Phenanthrene-d10	25.18	188	2443377	20.00	ng/uL	-0.06
59) Chrysene-d12	33.18	240	1512116	20.00	ng/uL	-0.06
68) Perylene-d12	37.27	264	1020421	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	0.00	99	0	0.00	ng/uL	
Spiked Amount 75.000	Range 19 - 32		Recovery =	0.00%#		
6) 2-Chlorophenol-d4	11.89	132	671	0.01	ng/uL	0.47
Spiked Amount 75.000	Range 34 - 84		Recovery =	0.01%#		
7) 1,2-Dichlorobenzene-d4	12.17	152	911	0.02	ng/uL	-0.29
Spiked Amount 50.000	Range 26 - 73		Recovery =	0.04%#		
20) Nitrobenzene-d5	0.00	82	0	0.00	ng/uL	
Spiked Amount 50.000	Range 55 - 98		Recovery =	0.00%#		
32) 2-Fluorobiphenyl	18.93	172	1763	0.02	ng/uL	0.10
Spiked Amount 50.000	Range 42 - 78		Recovery =	0.04%#		
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	731	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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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	0.00	178	0	N.D.	
56)	Anthracene	0.00	178	0	N.D.	
57)	Di-n-butylphthalate	27.18	149	30339	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	3100	N.D.	
66)	Bis(2-ethylhexyl)phthalate	33.45	149	27372	0.31 ng/uL	99
67)	Chrysene	33.18	228	3100	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
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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	37.27	252	2853	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.	

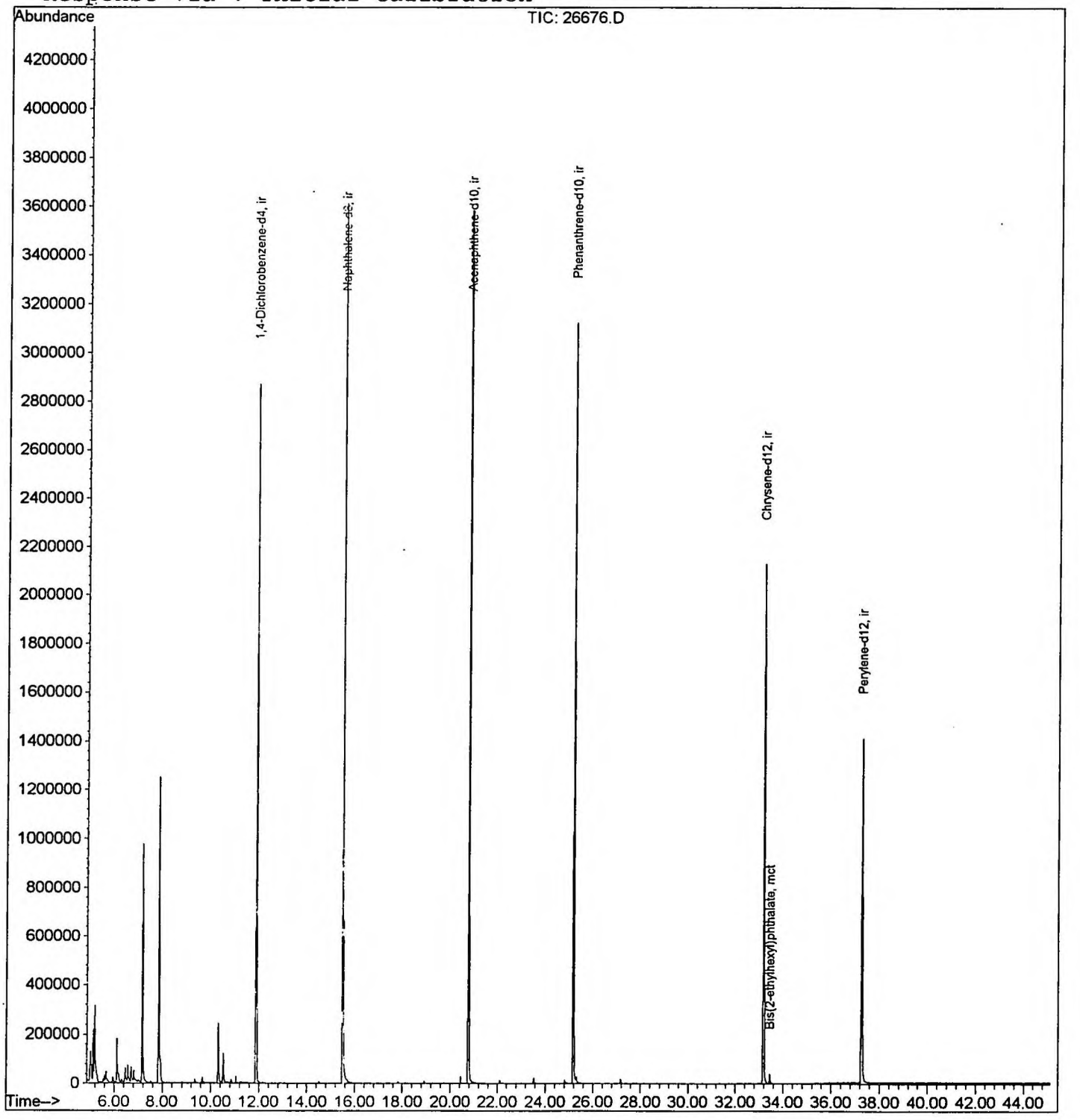
(#) = qualifier out of range (m) = manual integration

Quantitation Report

Data File : C:\HPCHEM\1\DATA\DEC0601\26676.D
Acq On : 6 Dec 2001 5:28 pm
Sample : gc/ms unknown
Misc :
MS Integration Params: RTEINT.P
Quant Time: Dec 7 10:14 2001

Vial: 9
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 : Fri Dec 07 09:59:59 2001
Response via : Initial Calibration



LSC Area Percent Report

Data File : C:\HPCHEM\1\DATA\DEC0601\26676.D Vial: 9
Acq On : 6 Dec 2001 5:28 pm Operator: GALOS
Sample : gc/ms unknown 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

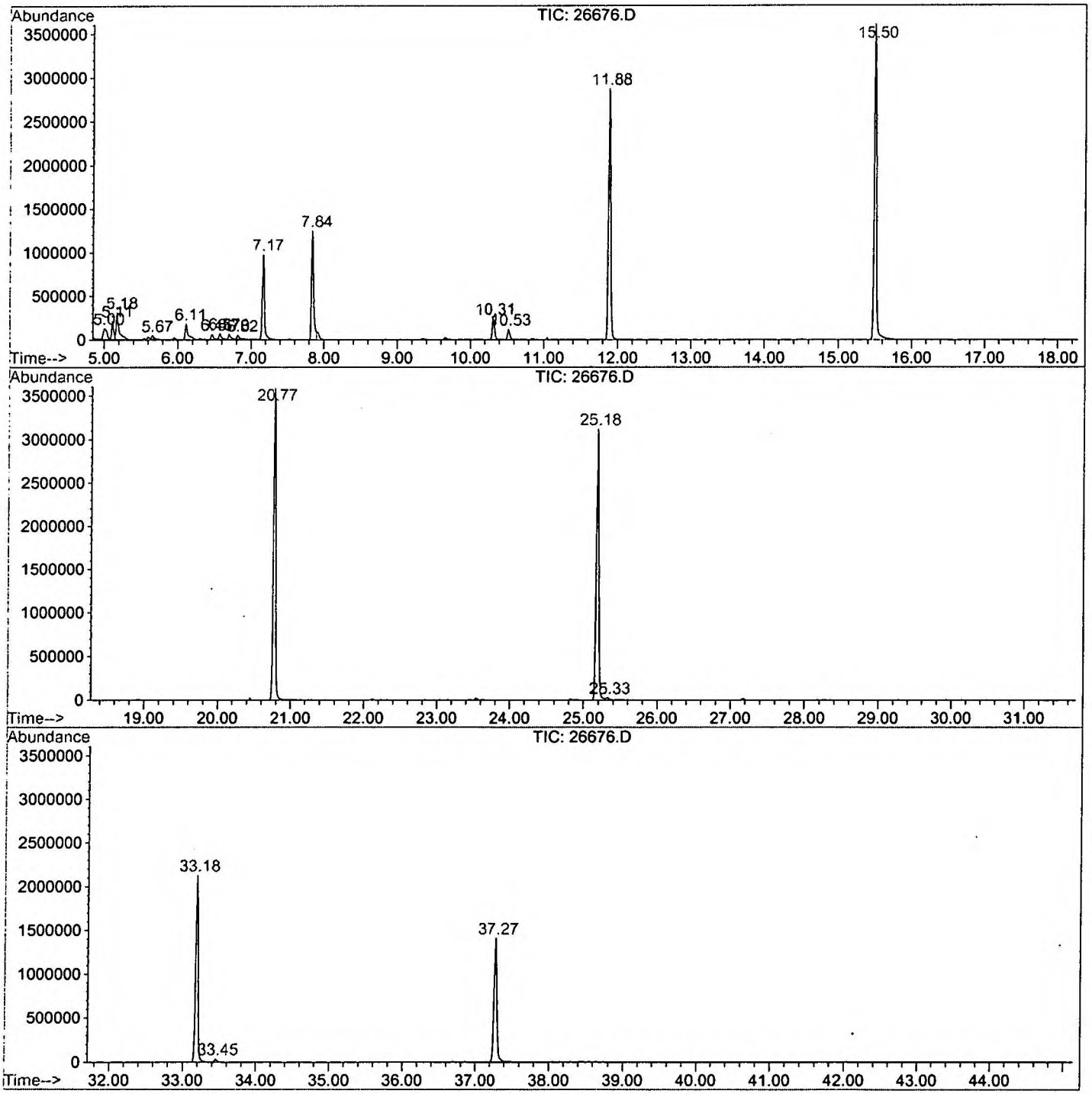
peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	peak area	peak % max.	% of total
1	5.005	12	18	27	rBV	123388	438200	5.75%	0.975%
2	5.115	27	30	34	rVV	210687	392402	5.15%	0.873%
3	5.179	34	37	55	rVB	310749	839262	11.01%	1.866%
4	5.665	86	90	95	rVB	40025	83207	1.09%	0.185%
5	6.115	135	139	156	rBV	182231	495781	6.50%	1.103%
6	6.463	174	177	186	rBV	58968	142481	1.87%	0.317%
7	6.573	186	189	200	rVB	68473	151416	1.99%	0.337%
8	6.701	200	203	211	rBV	60747	138267	1.81%	0.307%
9	6.821	212	216	220	rBV2	42997	79887	1.05%	0.178%
10	7.169	248	254	278	rVB	977976	1963120	25.75%	4.366%
11	7.839	323	327	334	rBV	1249675	2482114	32.56%	5.520%
12	10.314	593	597	609	rBV	243912	548910	7.20%	1.221%
13	10.525	616	620	628	rBV	121136	253594	3.33%	0.564%
14	11.883	763	768	782	rBV	2870032	6192722	81.23%	13.772%
15	15.496	1156	1162	1181	rBV	3611082	7624112	100.00%	16.955%
16	20.768	1731	1737	1755	rBV	3584557	7583398	99.47%	16.865%
17	25.179	2212	2218	2230	rBV2	3122134	6846138	89.80%	15.225%
18	25.326	2230	2234	2243	rVB2	27329	81024	1.06%	0.180%
19	33.185	3084	3091	3109	rBV2	2128609	4816467	63.17%	10.711%
20	33.451	3116	3120	3130	rVB	38487	90668	1.19%	0.202%
21	37.274	3529	3537	3550	rBV	1406787	3722299	48.82%	8.278%

Sum of corrected areas: 44965469

26676.D NP112701.M Fri Dec 07 14:51:16 2001

LSC Report - Integrated Chromatogram

File : C:\HPCHEM\1\DATA\DEC0601\26676.D
Operator : GALOS
Acquired : 6 Dec 2001 5:28 pm using AcqMethod NP112701
Instrument : GC/MS Ins
Sample Name: gc/ms unknown
Misc Info :
Vial Number: 9
Quant File :NP112701.RES (RTE Integrator)



26676.D NP112701.M Fri Dec 07 14:51:18 2001

Library Search Compound Report

Data File : C:\HPCHEM\1\DATA\DEC0601\26676.D
Acq On : 6 Dec 2001 5:28 pm
Sample : gc/ms unknown
Misc :
MS Integration Params: LSCINT.P

Vial: 9
Operator: GALOS
Inst : GC/MS Ins
Multiplr: 1.00

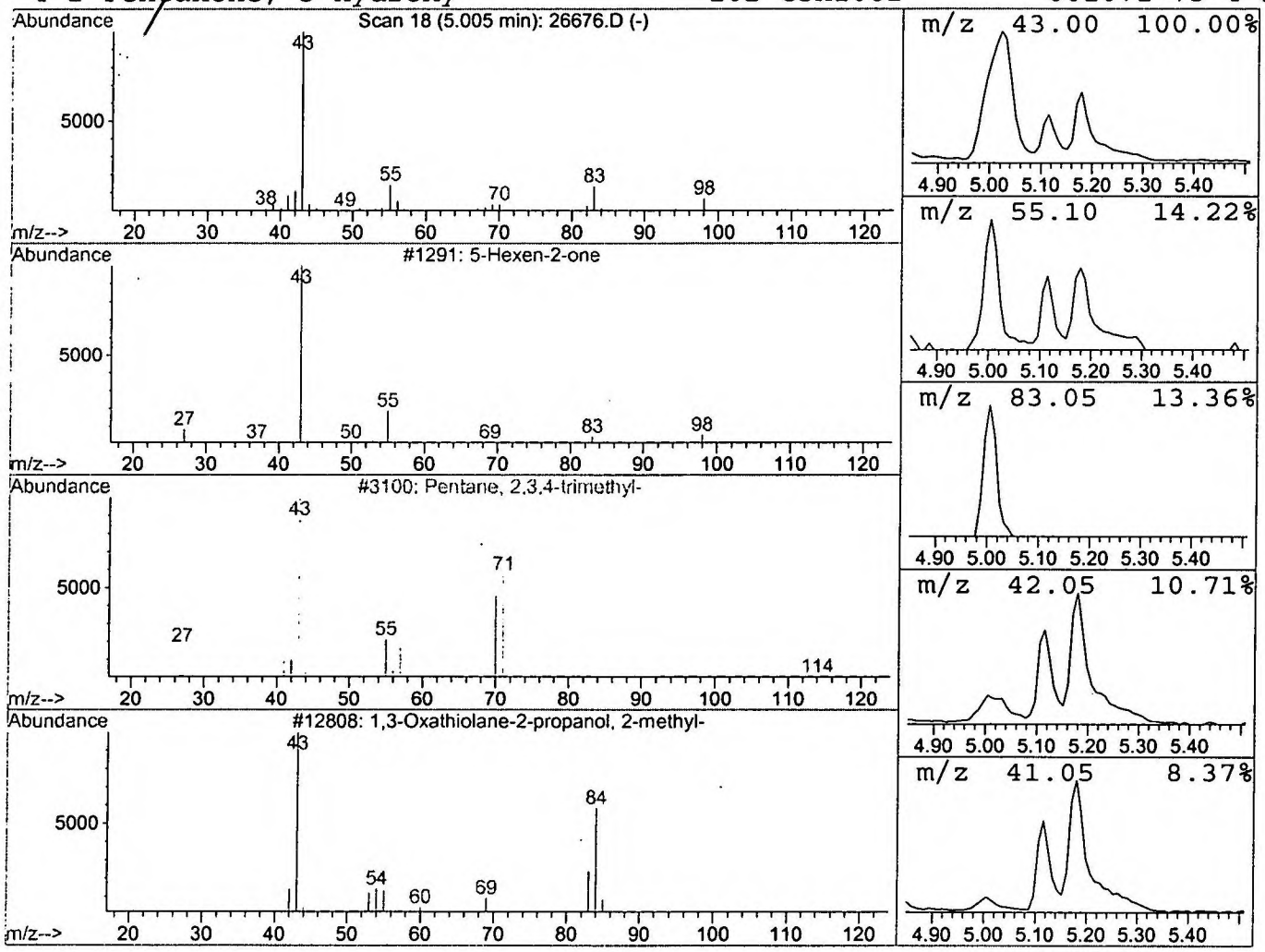
Quant Method : C:\HPCHEM\1\METHODS\NP112701.M (RTE Integrator)
Title : 208 625/TCLP calibration table
Library : C:\DATABASE\NBS75K.L

****

Peak Number 1 5-Hexen-2-one Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.00	1.42 ng/uL	438200	1,4-Dichlorobenzene-d4	11.89

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	5-Hexen-2-one	98	C6H10O	000109-49-9	9
2	Pentane, 2,3,4-trimethyl-	114	C8H18	000565-75-3	9
3	1,3-Oxathiolane-2-propanol, 2-methy	162	C7H14O2S	036651-30-6	9
4	2-Pentanone, 5-hydroxy-	102	C5H10O2	001071-73-4	9



Library Search Compound Report

Data File : C:\HPCHEM\1\DATA\DEC0601\26676.D
Acq On : 6 Dec 2001 5:28 pm
Sample : gc/ms unknown
Misc :
MS Integration Params: LSCINT.P

Vial: 9
Operator: GALOS
Inst : GC/MS Ins
Multiplr: 1.00

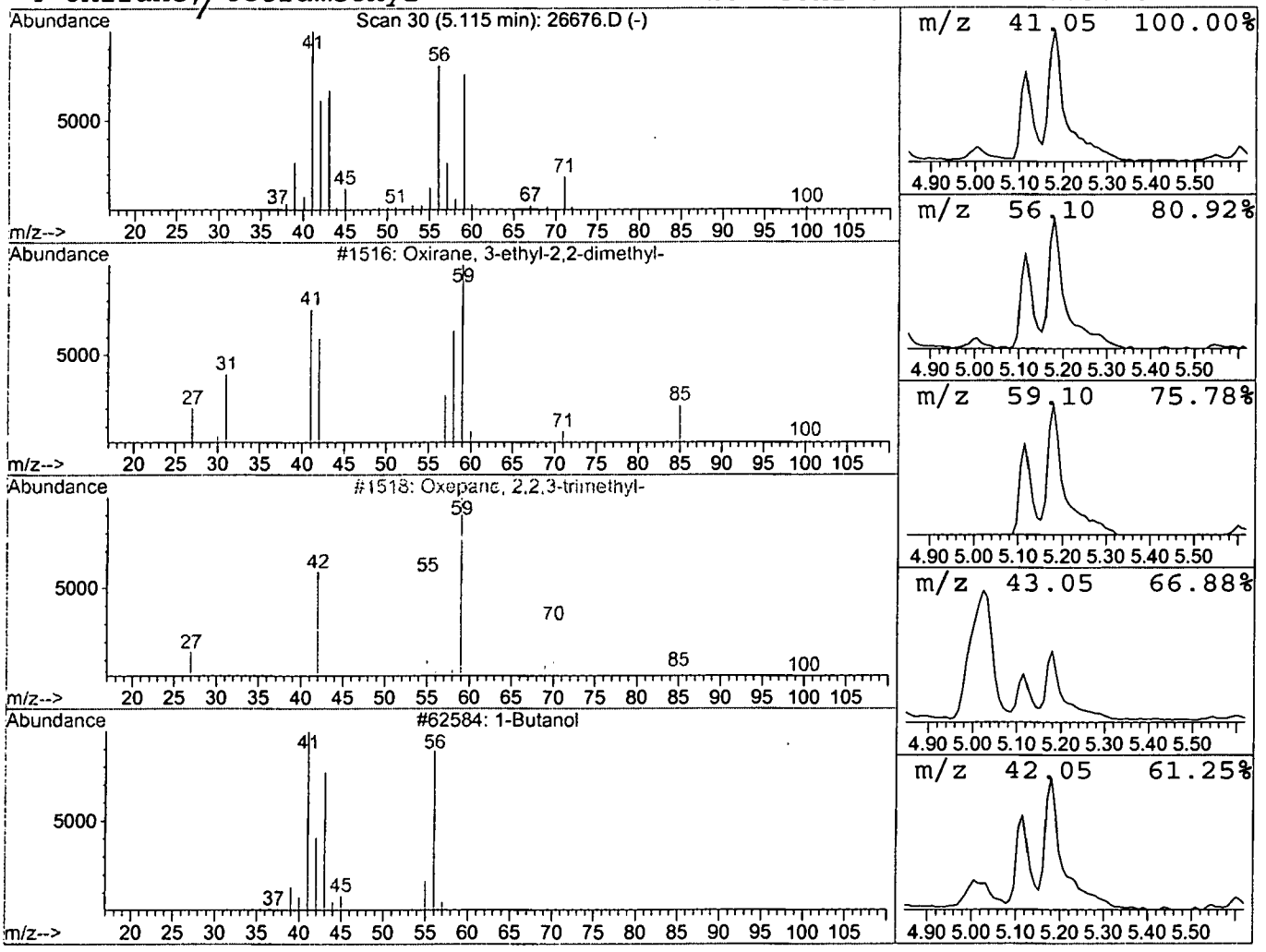
Quant Method : C:\HPCHEM\1\METHODS\NP112701.M (RTE Integrator)
Title : 208 625/TCLP calibration table
Library : C:\DATABASE\NBS75K.L

****

Peak Number 2 Oxirane, 3-ethyl-2,2-dimethyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.11	1.27 ng/uL	392402	1,4-Dichlorobenzene-d4	11.89

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Oxirane, 3-ethyl-2,2-dimethyl-	100	C6H12O	001192-22-9	43
2	Oxepane, 2,2,3-trimethyl-	100	C6H12O	023120-43-6	38
3	1-Butanol	74	C4H10O	000071-36-3	32
4	Oxirane, tetramethyl-	100	C6H12O	005076-20-0	27



Library Search Compound Report

Data File : C:\HPCHEM\1\DATA\DEC0601\26676.D
Acq On : 6 Dec 2001 5:28 pm
Sample : gc/ms unknown
Misc :
MS Integration Params: LSCINT.P

Vial: 9
Operator: GALOS
Inst : GC/MS Ins
Multiplr: 1.00

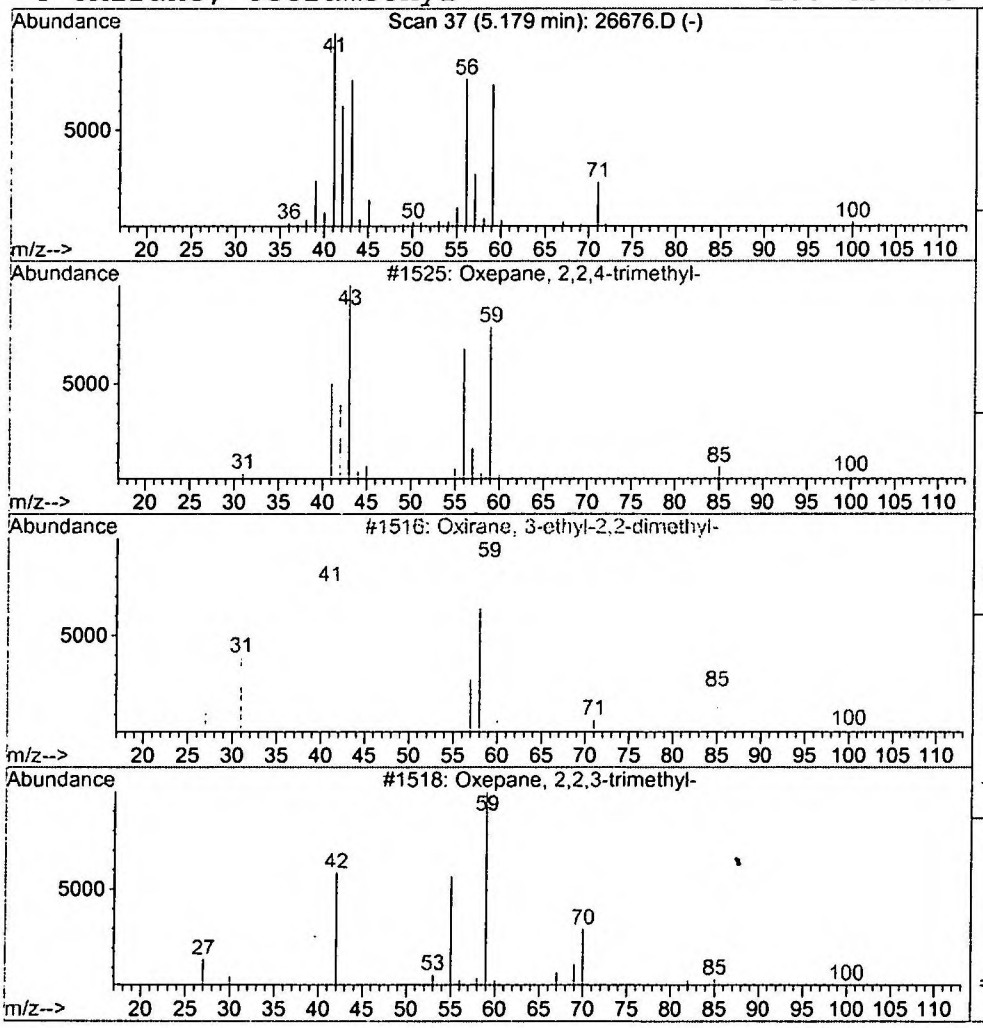
Quant Method : C:\HPCHEM\1\METHODS\NP112701.M (RTE Integrator)
Title : 208 625/TCLP calibration table
Library : C:\DATABASE\NBS75K.L

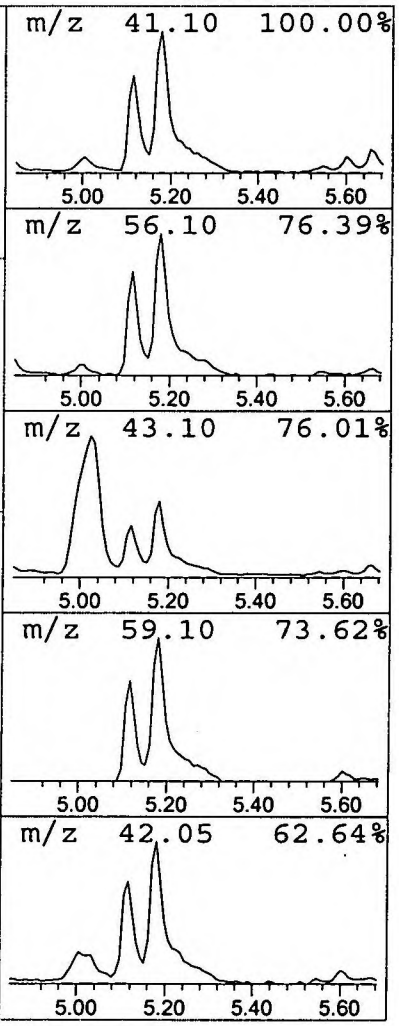
****

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.18	2.71 ng/uL	839262	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	Oxepane, 2,2,3-trimethyl-	100	C6H12O	023120-43-6	35
4	Oxirane, tetramethyl-	100	C6H12O	005076-20-0	35





Library Search Compound Report

Data File : C:\HPCHEM\1\DATA\DEC0601\26676.D
Acq On : 6 Dec 2001 5:28 pm
Sample : gc/ms unknown
Misc :
MS Integration Params: LSCINT.P

Vial: 9
Operator: GALOS
Inst : GC/MS Ins
Multiplr: 1.00

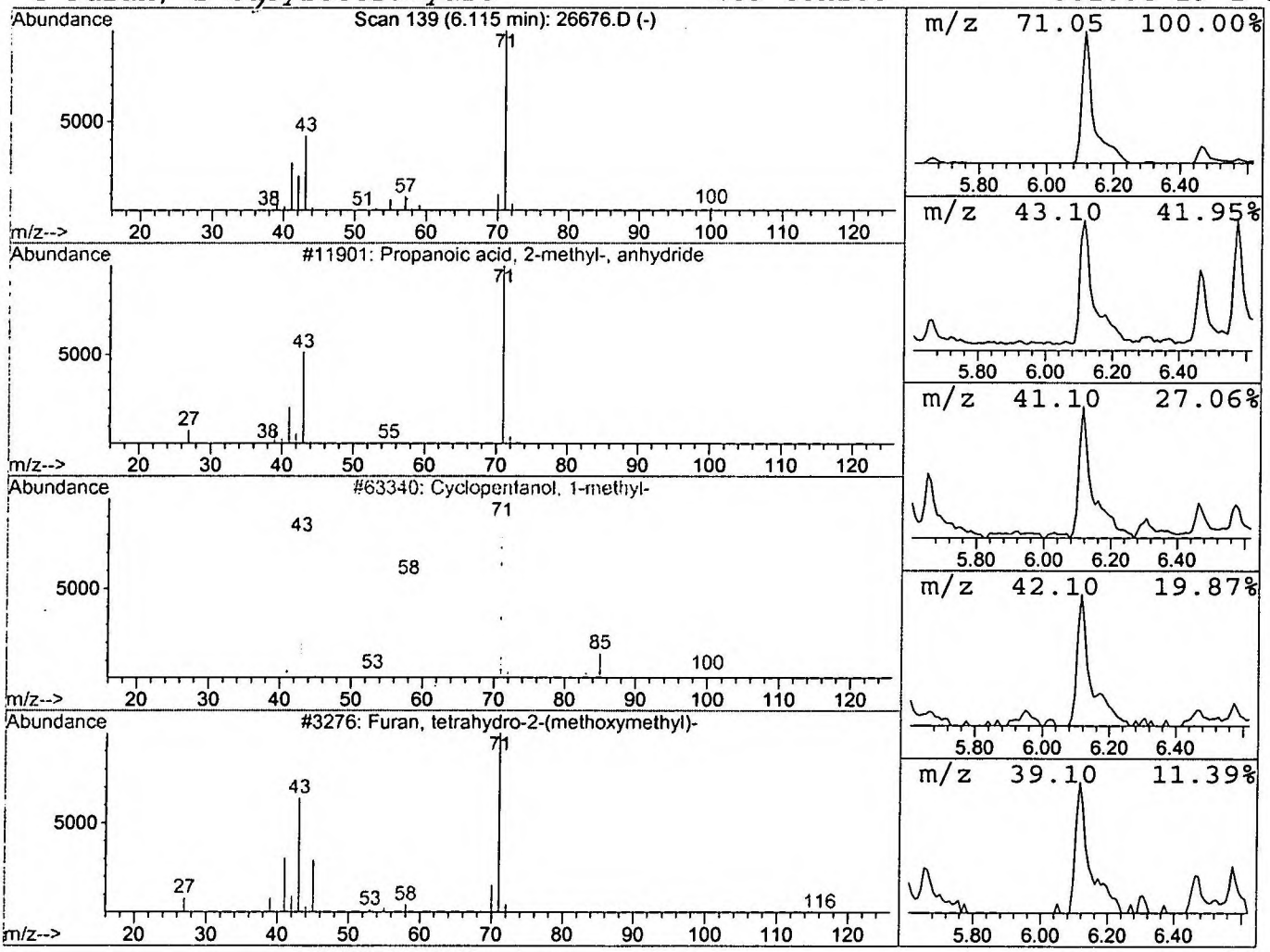
Quant Method : C:\HPCHEM\1\METHODS\NP112701.M (RTE Integrator)
Title : 208 625/TCLP calibration table
Library : C:\DATABASE\NBS75K.L

****

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.11	1.60 ng/uL	495781	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	42
2	Cyclopentanol, 1-methyl-	100	C6H12O	001462-03-9	35
3	Furan, tetrahydro-2-(methoxymethyl)	116	C6H12O2	019354-27-9	9
4	Furan, 2-butyltetrahydro-	128	C8H16O	001004-29-1	9



Library Search Compound Report

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

Vial: 9
Operator: GALOS
Inst : GC/MS Ins
Multiplr: 1.00

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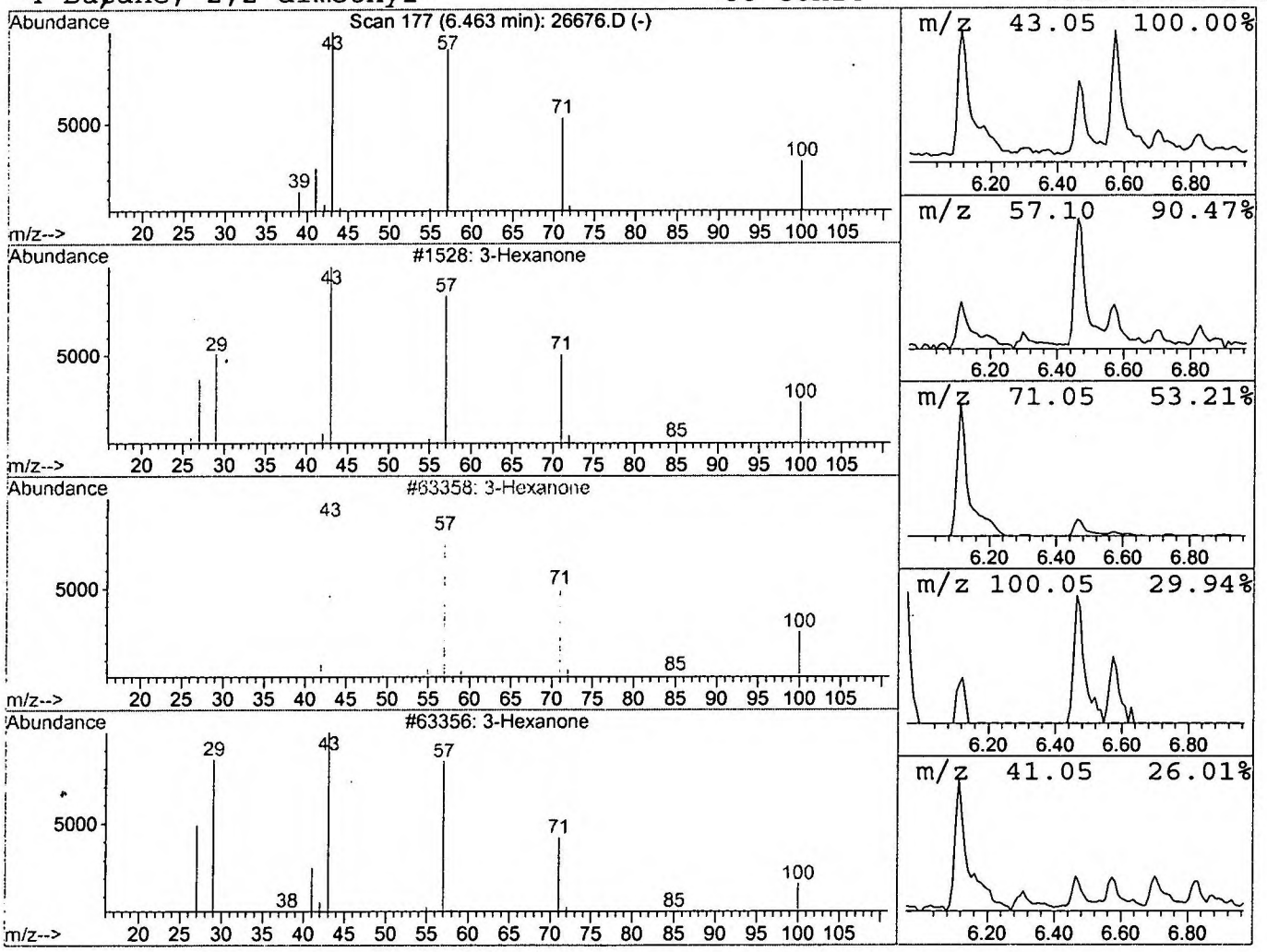
Library : C:\DATABASE\NBS75K.L

****

Peak Number 5 3-Hexanone Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.46	0.46 ng/uL	142481	1,4-Dichlorobenzene-d4	11.89

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	3-Hexanone	100	C6H12O	000589-38-8	90
2	3-Hexanone	100	C6H12O	000589-38-8	90
3	3-Hexanone	100	C6H12O	000589-38-8	72
4	Butane, 2,2-dimethyl-	86	C6H14	000075-83-2	59



Library Search Compound Report

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Acq On : 6 Dec 2001 5:28 pm Operator: GALOS
Sample : gc/ms unknown Inst : GC/MS Ins
Misc : Multiplr: 1.00
MS Integration Params: LSCINT.P

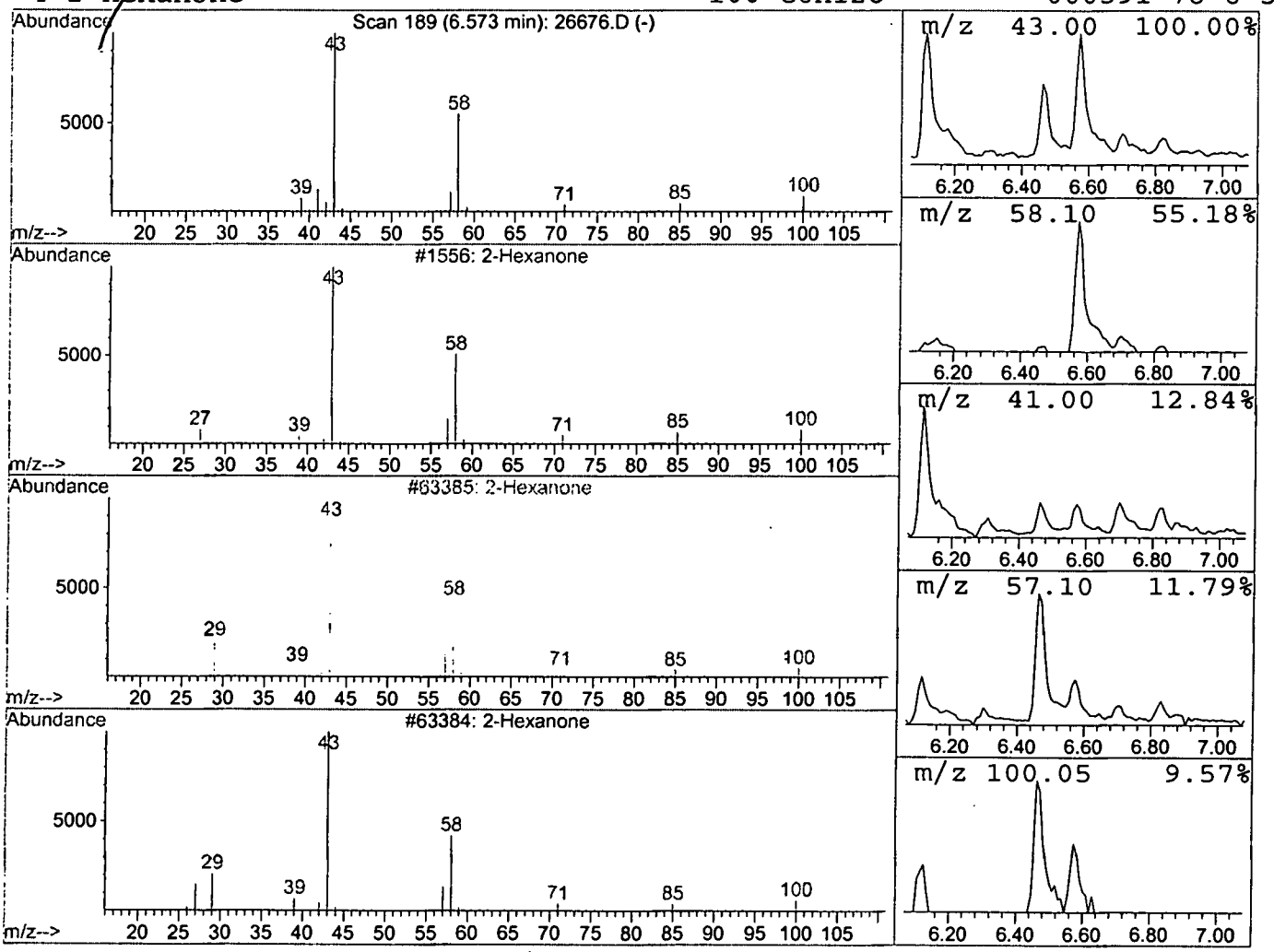
Quant Method : C:\HPCHEM\1\METHODS\NP112701.M (RTE Integrator)
Title : 208 625/TCLP calibration table
Library : C:\DATABASE\NBS75K.L

****

Peak Number 6 2-Hexanone Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.57	0.49 ng/uL	151416	1,4-Dichlorobenzene-d4	11.89

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Hexanone	100	C6H12O	000591-78-6	72
2	2-Hexanone	100	C6H12O	000591-78-6	64
3	2-Hexanone	100	C6H12O	000591-78-6	64
4	2-Hexanone	100	C6H12O	000591-78-6	59



Library Search Compound Report

Data File : C:\HPCHEM\1\DATA\DEC0601\26676.D Vial: 9
Acq On : 6 Dec 2001 5:28 pm Operator: GALOS
Sample : gc/ms unknown Inst : GC/MS Ins
Misc : Multiplr: 1.00
MS Integration Params: LSCINT.P

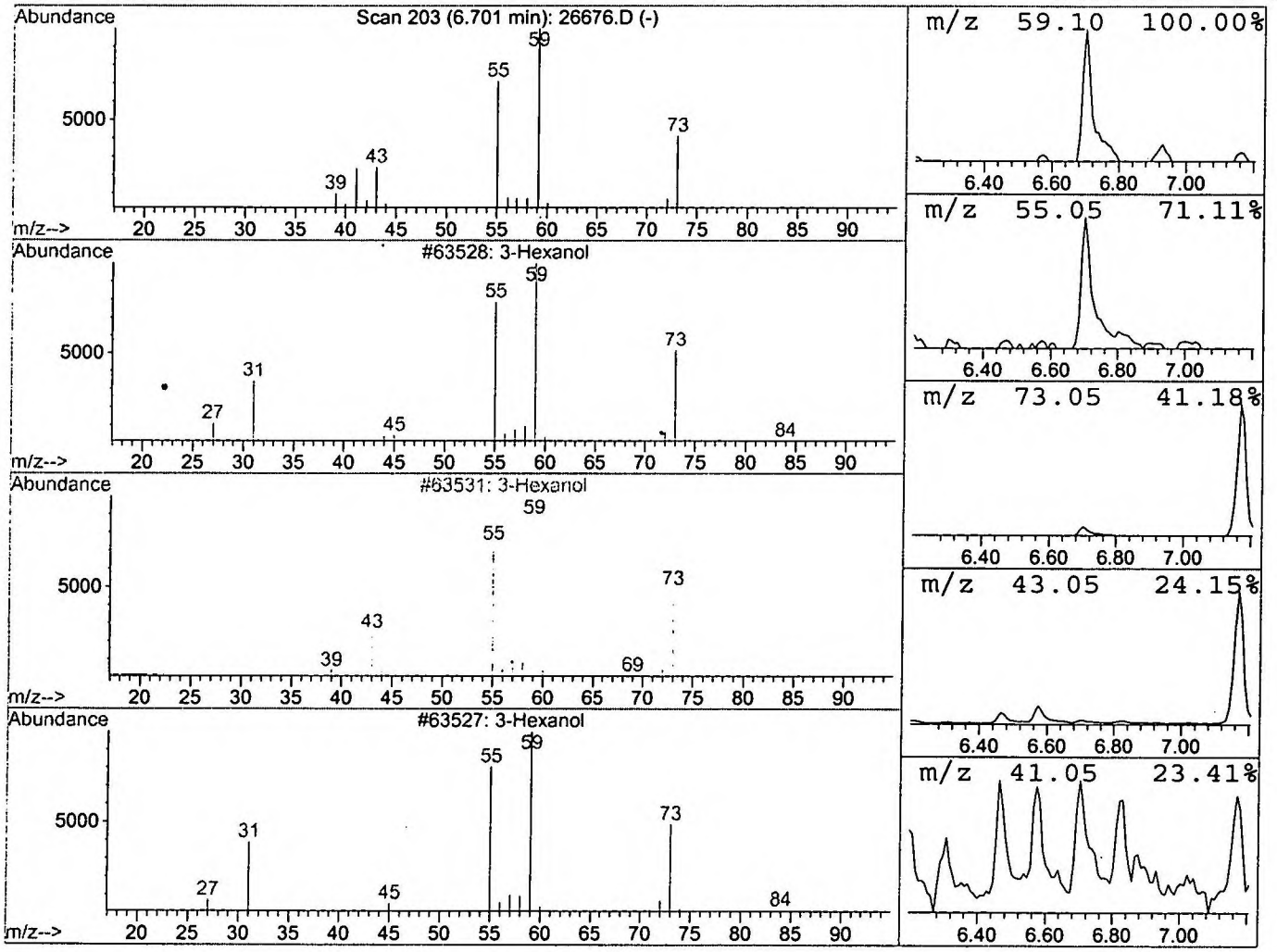
Quant Method : C:\HPCHEM\1\METHODS\NP112701.M (RTE Integrator)
Title : 208 625/TCLP calibration table
Library : C:\DATABASE\NBS75K.L

****

Peak Number 7 3-Hexanol Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.70	0.45 ng/uL	138267	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	64
4	3-Hexanol	102	C6H14O	000623-37-0	64



Library Search Compound Report

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Acq On : 6 Dec 2001 5:28 pm
Sample : gc/ms unknown
Misc :
MS Integration Params: LSCINT.P

Vial: 9
Operator: GALOS
Inst : GC/MS Ins
Multiplr: 1.00

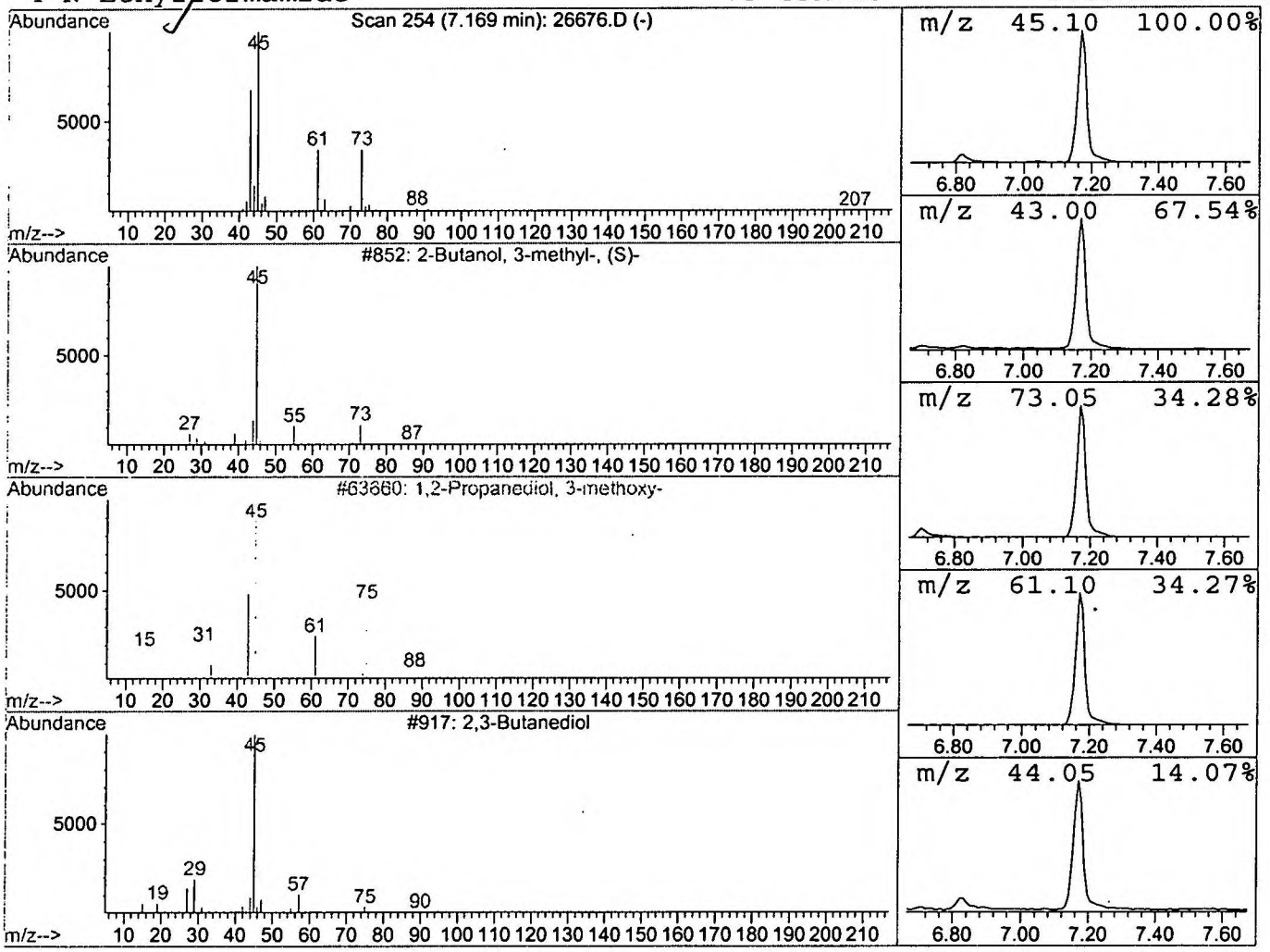
Quant Method : C:\HPCHEM\1\METHODS\NP112701.M (RTE Integrator)
Title : 208 625/TCLP calibration table
Library : C:\DATABASE\NBS75K.L

****

Peak Number 8 2-Butanol, 3-methyl-, (S)- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.17	6.34 ng/uL	1963120	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	1,2-Propanediol, 3-methoxy-	106	C4H10O3	000623-39-2	36
3	2,3-Butanediol	90	C4H10O2	000513-85-9	33
4	N-Ethylformamide	73	C3H7NO	000627-45-2	9



Library Search Compound Report

Data File : C:\HPCHEM\1\DATA\DEC0601\26676.D Vial: 9
Acq On : 6 Dec 2001 5:28 pm Operator: GALOS
Sample : gc/ms unknown Inst : GC/MS Ins
Misc : Multiplr: 1.00
MS Integration Params: LSCINT.P

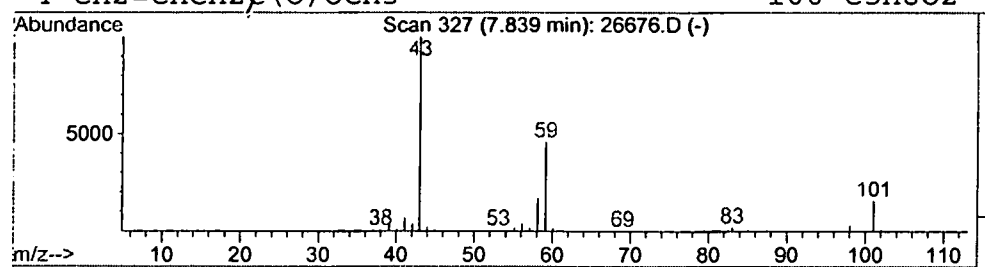
Quant Method : C:\HPCHEM\1\METHODS\NP112701.M (RTE Integrator)
Title : 208 625/TCLP calibration table
Library : C:\DATABASE\NBS75K.L

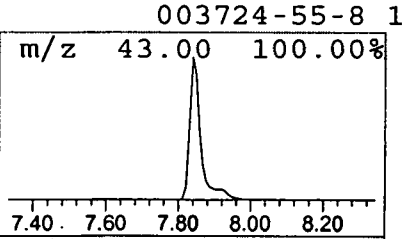
****

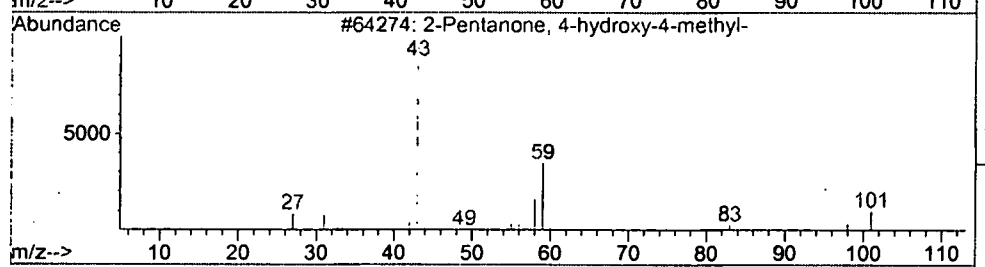
Peak Number 9 2-Pentanone, 4-hydroxy-4-methy Concentration Rank 1

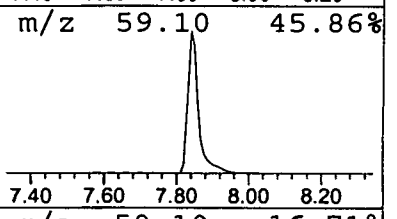
R.T.	EstConc	Area	Relative to ISTD	R.T.
7.84	8.02 ng/uL	2482110	1,4-Dichlorobenzene-d4	11.89

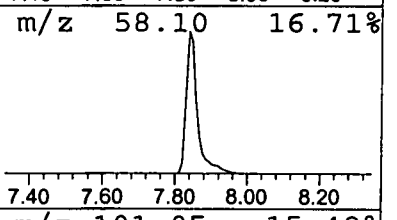
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	64
2	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	43
3	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	25
4	CH2=CHCH2C(O)OCH3	100	C5H8O2	003724-55-8	10

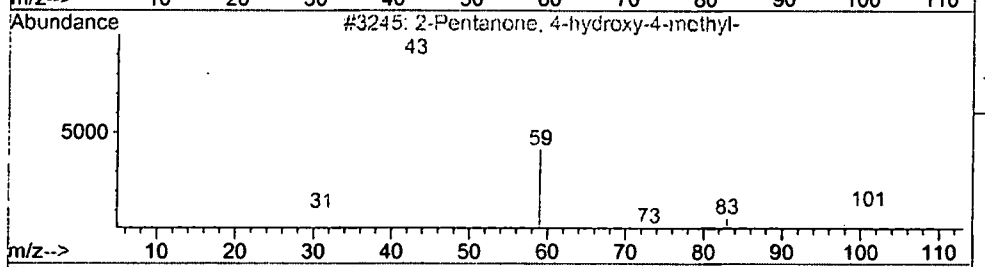


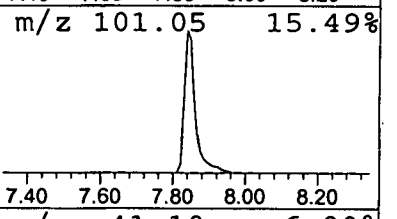


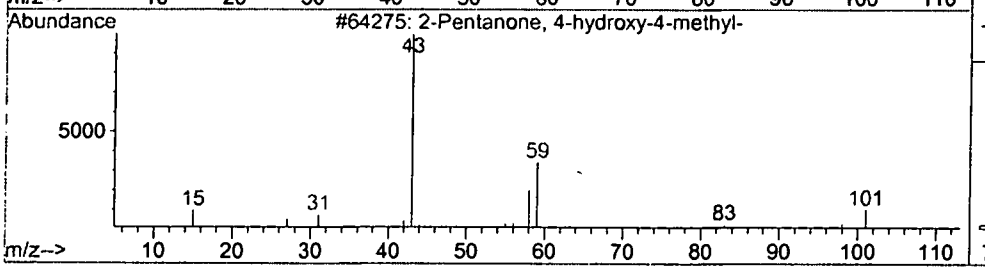


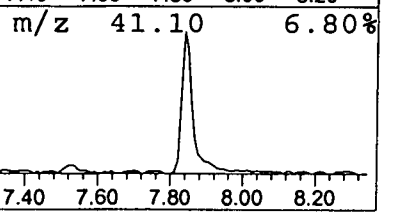












Library Search Compound Report

Data File : C:\HPCHEM\1\DATA\DEC0601\26676.D
Acq On : 6 Dec 2001 5:28 pm
Sample : gc/ms unknown
Misc :
MS Integration Params: LSCINT.P

Vial: 9
Operator: GALOS
Inst : GC/MS Ins
Multiplr: 1.00

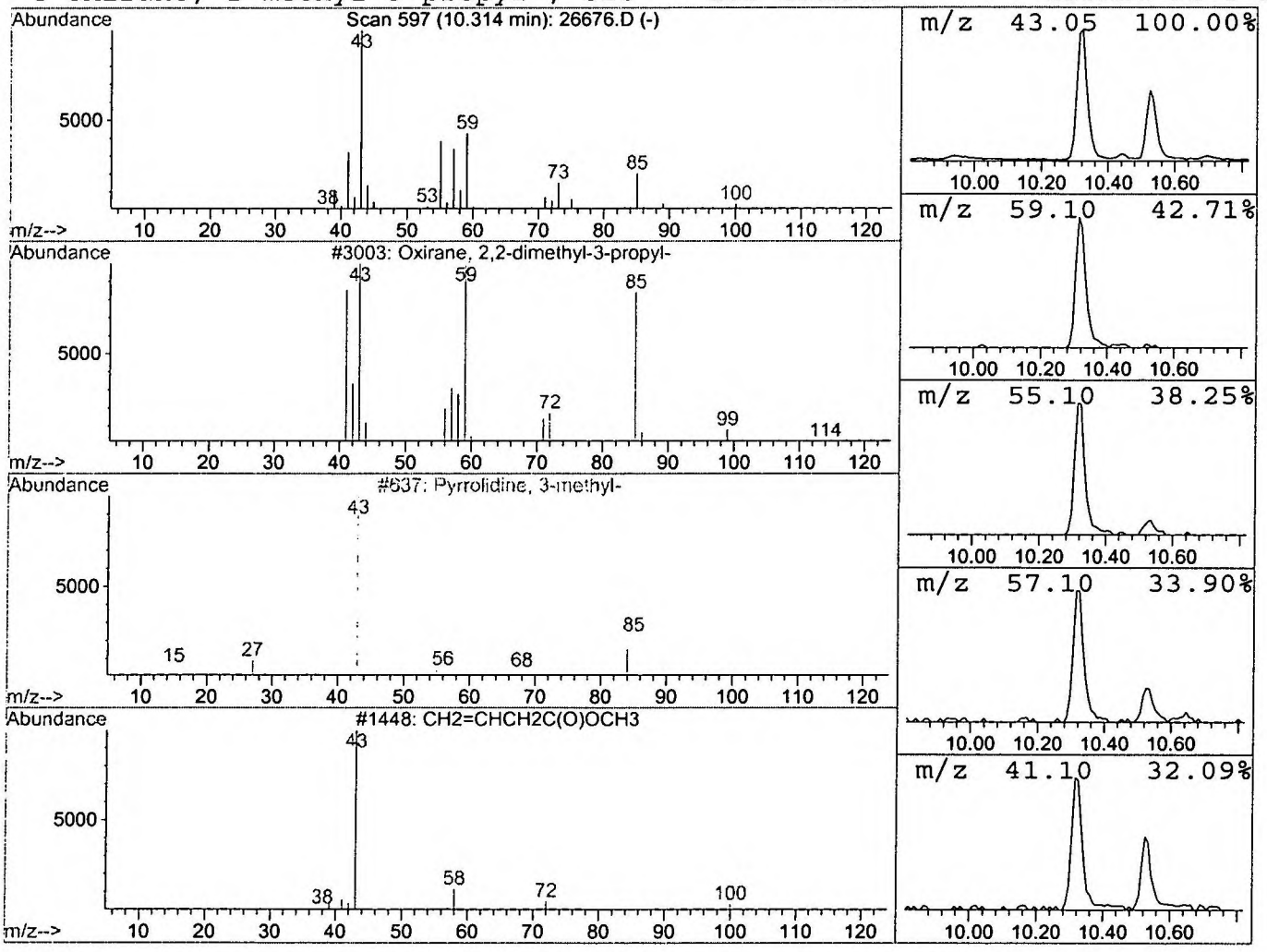
Quant Method : C:\HPCHEM\1\METHODS\NP112701.M (RTE Integrator)
Title : 208 625/TCLP calibration table
Library : C:\DATABASE\NBS75K.L

****

Peak Number 10 Oxirane, 2,2-dimethyl-3-propyl Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.31	1.77 ng/uL	548910	1,4-Dichlorobenzene-d4	11.89

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Oxirane, 2,2-dimethyl-3-propyl-	114	C7H14O	017612-35-0	17
2	Pyrrolidine, 3-methyl-	85	C5H11N	034375-89-8	14
3	CH2=CHCH2C(O)OCH3	100	C5H8O2	003724-55-8	14
4	Oxirane, 2-methyl-3-propyl-, cis-	100	C6H12O	006124-90-9	12



Library Search Compound Report

Data File : C:\HPCHEM\1\DATA\DEC0601\26676.D Vial: 9
Acq On : 6 Dec 2001 5:28 pm Operator: GALOS
Sample : gc/ms unknown Inst : GC/MS Ins
Misc : Multiplr: 1.00
MS Integration Params: LSCINT.P

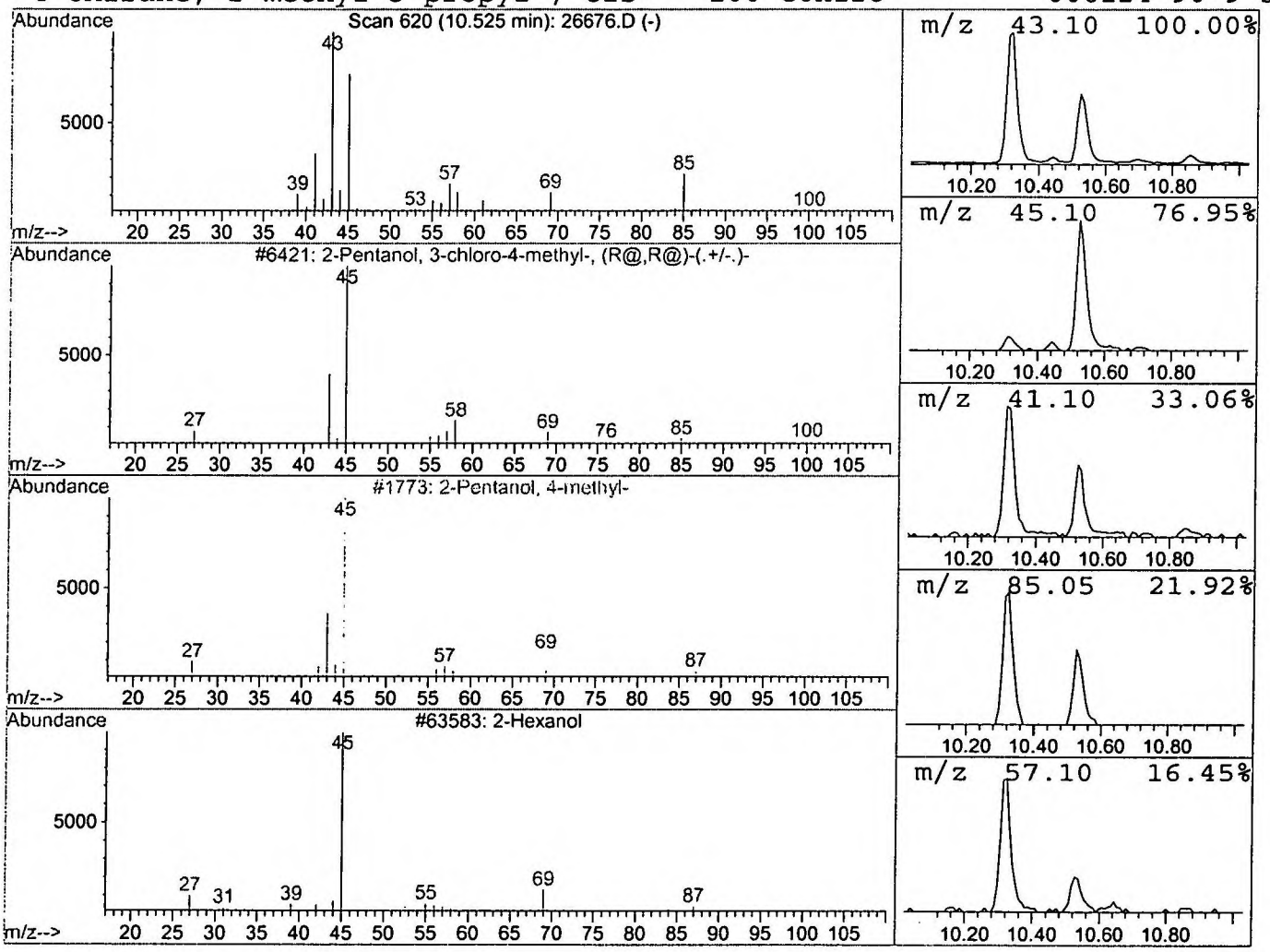
Quant Method : C:\HPCHEM\1\METHODS\NP112701.M (RTE Integrator)
Title : 208 625/TCLP calibration table
Library : C:\DATABASE\NBS75K.L

****

Peak Number 11 2-Pentanol, 3-chloro-4-methyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.53	0.82 ng/uL	253594	1,4-Dichlorobenzene-d4	11.89

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Pentanol, 3-chloro-4-methyl-, (R@	136	C6H13ClO	074685-47-5	53
2	2-Pentanol, 4-methyl-	102	C6H14O	000108-11-2	40
3	2-Hexanol	102	C6H14O	000626-93-7	38
4	Oxirane, 2-methyl-3-propyl-, cis-	100	C6H12O	006124-90-9	37



Tentatively Identified Compound (LSC) summary

Operator ID: GALOS Date Acquired: 6 Dec 2001 5:28 pm
Data File: C:\HPCHEM\1\DATA\DEC0601\26676.D
Name: gc/ms unknown
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
5-Hexen-2-one	5.00	1.4	ng/uL	438200	ISTD01	11.89	6192720	20.0
Oxirane, 3-ethyl-2,2	5.11	1.3	ng/uL	392402	ISTD01	11.89	6192720	20.0
Oxepane, 2,2,4-trime	5.18	2.7	ng/uL	839262	ISTD01	11.89	6192720	20.0
Propanoic acid, 2-me	6.11	1.6	ng/uL	495781	ISTD01	11.89	6192720	20.0
3-Hexanone	6.46	0.5	ng/uL	142481	ISTD01	11.89	6192720	20.0
2-Hexanone	6.57	0.5	ng/uL	151416	ISTD01	11.89	6192720	20.0
3-Hexanol	6.70	0.4	ng/uL	138267	ISTD01	11.89	6192720	20.0
2-Butanol, 3-methyl-	7.17	6.3	ng/uL	1963120	ISTD01	11.89	6192720	20.0
2-Pentanone, 4-hydro	7.84	8.0	ng/uL	2482110	ISTD01	11.89	6192720	20.0
Oxirane, 2,2-dimethy	10.31	1.8	ng/uL	548910	ISTD01	11.89	6192720	20.0
2-Pentanol, 3-chloro	10.53	0.8	ng/uL	253594	ISTD01	11.89	6192720	20.0

26676.D NP112701.M Fri Dec 07 14:51:42 2001

Comments for Sample AD26676 - Analysis $GCMS

Westchester Dept. of Labs & Research
User: Vinci, David L
Date: 12-16-2001 Time: 11:25:12

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, except for the presence of bis(2-Ethylhexyl)phthalate at an estimated level of 0.31 ug/L.

0.33 in blank