Comments for Sample AD26475 - Analysis $GCMS

Westchester Dept. of Labs & Research
User: Vinci, David L
Date: 12-16-2001 Time: 08:32:44

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

User: Vinci, David L
Westchester Dept. of Labs & Research
Date: 12/16/2001 Time: 08:33:14

Sample: AD26475
Analysis: $GCMS GCMS Scan For Unknowns
Units: ug
Started: 12/16/01 08:30 Ended: 12/16/01 08:30 Analyst: DLV
Result source: manual entry
Page: 1

	Component Name	Viol	Result	Qualifier	MDL
1	Other unknown compounds		.		
2	Surr: 2,4-DDD	LSPEC	< MDL		5.0

Quantitation Report (QT Reviewed)

Data File : C:\HPCHEM\1\DATA\DEC0701\26475.D Vial: 5
Acq On : 7 Dec 2001 9:45 pm Operator: GALOS
Sample : gc/ms unknown 11/6 Inst : GC/MS Ins
Misc : Multiplr: 1.00
MS Integration Params: RTEINT.P
Quant Time: Dec 10 9:26 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	11.88	152	910895	20.00	ng/uL	-0.05
19) Naphthalene-d8	15.49	136	3480953	20.00	ng/uL	-0.05
31) Acenaphthene-d10	20.77	164	1590398	20.00	ng/uL	-0.05
49) Phenanthrene-d10	25.18	188	2333593	20.00	ng/uL	-0.06
59) Chrysene-d12	33.18	240	1564074	20.00	ng/uL	-0.06
68) Perylene-d12	37.26	264	1074965	20.00	ng/uL	-0.07
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	11.87	132	336	0.01	ng/uL	0.45
Spiked Amount 75.000	Range 34 - 84		Recovery =	0.01%#		
7) 1,2-Dichlorobenzene-d4	0.00	152	0	0.00	ng/uL	
Spiked Amount 50.000	Range 26 - 73		Recovery =	0.00%#		
20) Nitrobenzene-d5	0.00	82	0	0.00	ng/uL	
Spiked Amount 50.000	Range 55 - 98		Recovery =	0.00%#		
32) 2-Fluorobiphenyl	18.91	172	1655	0.02	ng/uL	0.08
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	12.88	45	1821	N.D.		
16) 3&4-Methylphenol	0.00	108	0	N.D.		

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

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	Compound	R.T.	QIon	Response	Conc Unit	Qvalue
17)	N-Nitrosodi-n-propylamine	12.72	70	1066	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.	
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	27.17	149	2482	N.D.	
58)	Fluoranthene	0.00	202	0	N.D.	
60)	3,3'-Dichlorobenzidine	0.00	252	0	N.D.	

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DataAcq Meth : NP112701

Compound	R.T.	QIon	Response	Conc Unit	Qvalue
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	33.17	228	3130	N.D.	
66) Bis(2-ethylhexyl)phthalate	33.43	149	1819	N.D.	
67) Chrysene	33.17	228	3130	N.D.	
69) Di-n-Octylphthalate	0.00	149	0	N.D.	
70) Benzo(b)fluoranthene	0.00	252	0	N.D.	
71) Benzo(k)fluoranthene	0.00	252	0	N.D.	
72) Benzo(a)pyrene	37.27	252	3266	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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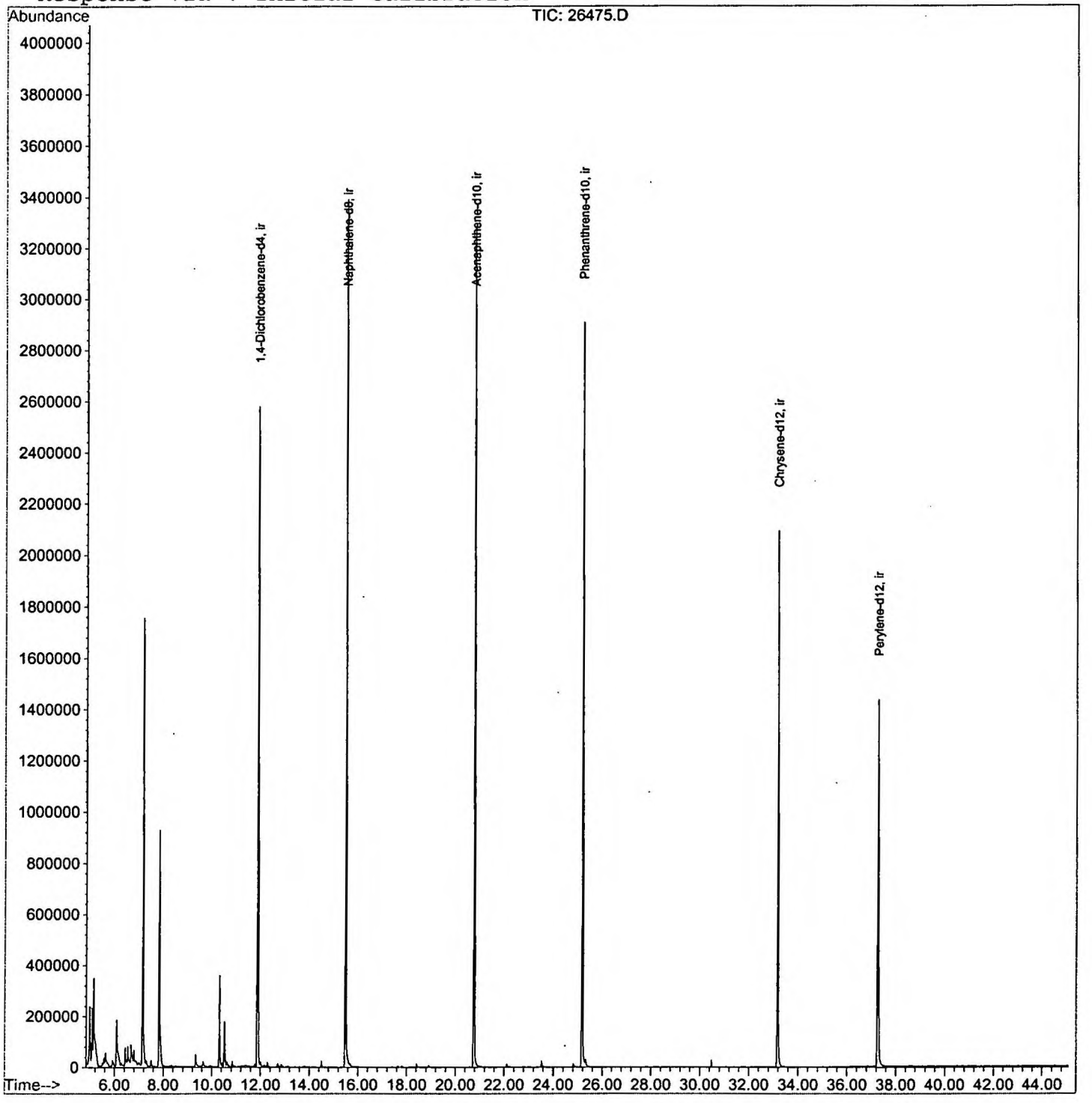
Misc :
MS Integration Params: RTEINT.P
Quant Time: Dec 10 9:26 2001

Vial: 5
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\DEC0701\26475.D Vial: 5
Acq On : 7 Dec 2001 9:45 pm Operator: GALOS
Sample : gc/ms unknown 11/6 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

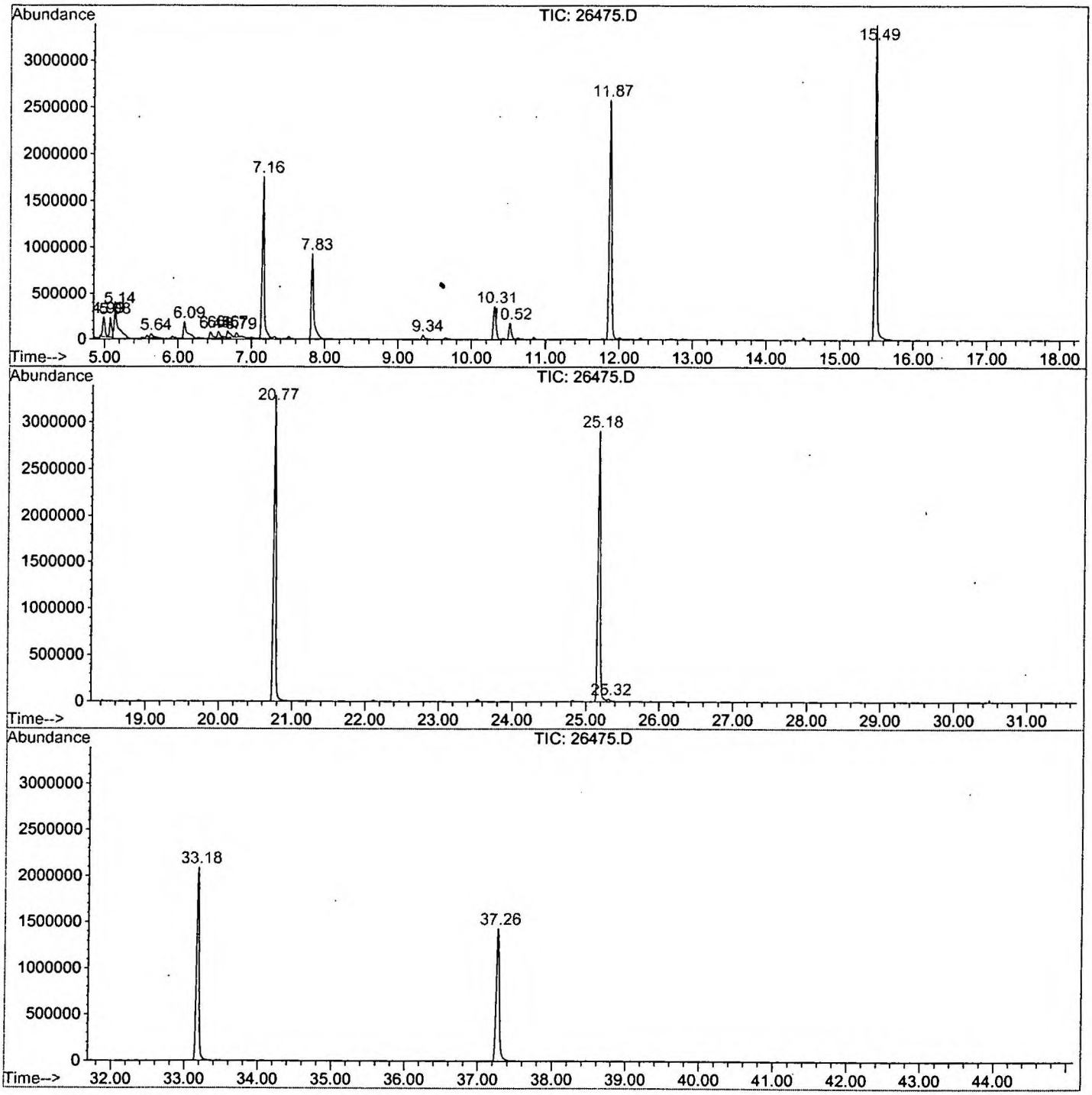
peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	peak area	peak % max.	% of total
1	4.985	12	16	23	rBV	215400	456682	6.49%	1.000%
2	5.077	23	26	30	rVV	215256	417216	5.93%	0.914%
3	5.141	30	33	56	rVB	345606	1242155	17.65%	2.721%
4	5.637	83	87	96	rVB3	44649	133131	1.89%	0.292%
5	6.086	132	136	154	rBV	181617	615331	8.74%	1.348%
6	6.444	171	175	182	rBV	71993	187057	2.66%	0.410%
7	6.554	183	187	197	rVV	70913	197213	2.80%	0.432%
8	6.673	197	200	210	rVV2	72879	235131	3.34%	0.515%
9	6.792	210	213	220	rVB2	45560	102676	1.46%	0.225%
10	7.159	247	253	268	rBV	1748880	3427834	48.70%	7.507%
11	7.828	322	326	343	rVB	925893	2113189	30.02%	4.628%
12	9.341	486	491	504	rBV	48619	124989	1.78%	0.274%
13	10.313	592	597	607	rBV	360015	822222	11.68%	1.801%
14	10.515	615	619	627	rBV	176473	396501	5.63%	0.868%
15	11.872	762	767	779	rBV	2578531	5614082	79.76%	12.296%
16	15.494	1156	1162	1175	rBV	3390599	7029157	99.86%	15.395%
17	20.767	1730	1737	1753	rBV	3284560	7038951	100.00%	15.416%
18	25.178	2211	2218	2230	rBV2	2907931	6538698	92.89%	14.321%
19	25.316	2230	2233	2245	rVB	29517	83905	1.19%	0.184%
20	33.183	3084	3091	3111	rBV2	2091949	4950688	70.33%	10.843%
21	37.264	3528	3536	3559	rBV2	1432008	3932084	55.86%	8.612%

Sum of corrected areas: 45658892

26475.D NP112701.M Thu Dec 13 11:19:18 2001

LSC Report - Integrated Chromatogram

File : C:\HPCHEM\1\DATA\DEC0701\26475.D
Operator : GALOS
Acquired : 7 Dec 2001 9:45 pm using AcqMethod NP112701
Instrument : GC/MS Ins
Sample Name: gc/ms unknown 11/6
Misc Info :
Vial Number: 5
Quant File :NP112701.RES (RTE Integrator)



26475.D NP112701.M Thu Dec 13 11:19:20 2001

Library Search Compound Report

Data File : C:\HPCHEM\1\DATA\DEC0701\26475.D
Acq On : 7 Dec 2001 9:45 pm
Sample : gc/ms unknown 11/6
Misc :
MS Integration Params: LSCINT.P

Vial: 5
Operator: GALOS
Inst : GC/MS Ins
Multiplr: 1.00

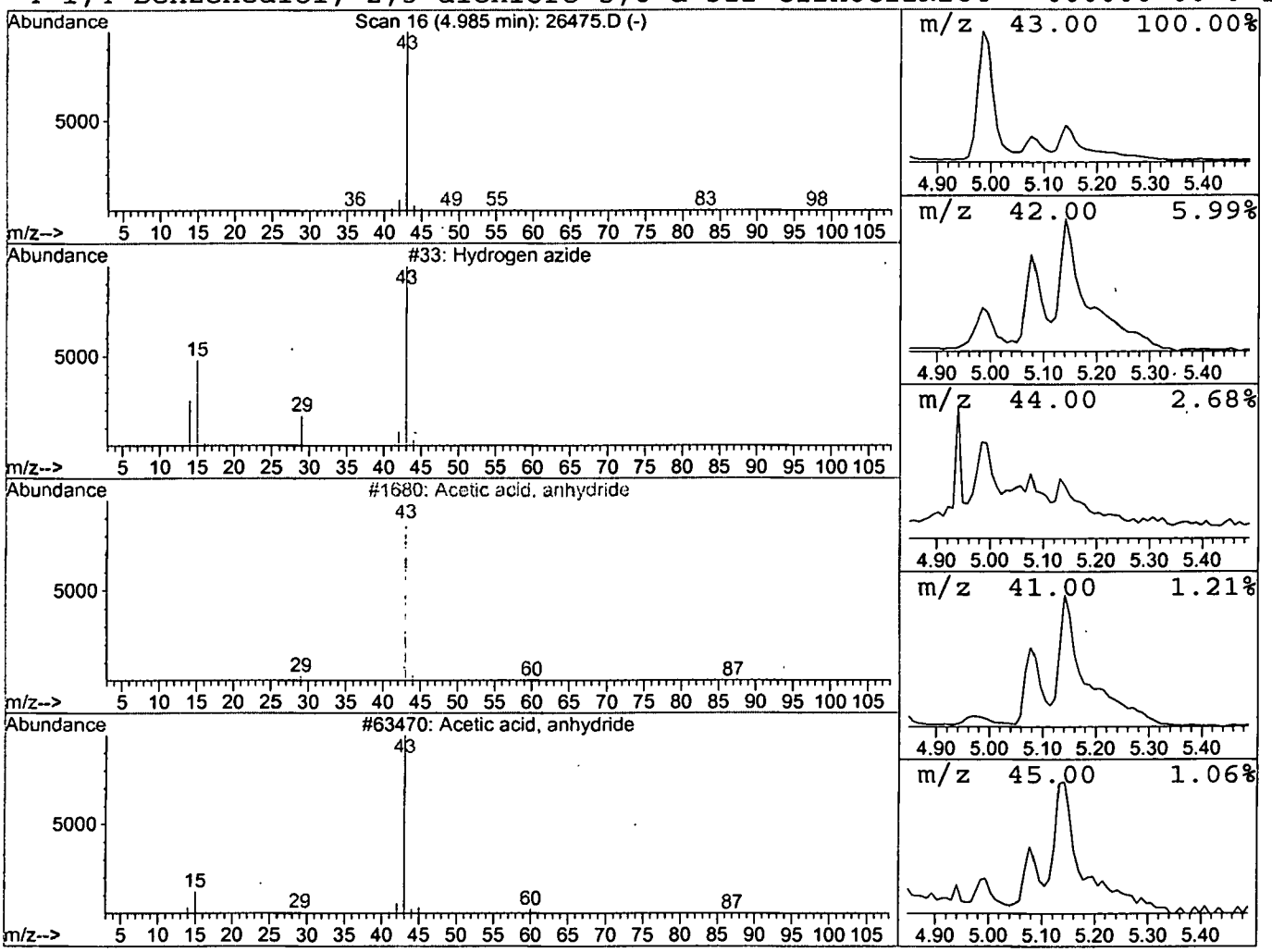
Quant Method : C:\HPCHEM\1\METHODS\NP112701.M (RTE Integrator)
Title : 208 625/TCLP calibration table
Library : C:\DATABASE\NBS75K.L

****

Peak Number 1 Hydrogen azide Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.99	1.63 ng/uL	456682	1,4-Dichlorobenzene-d4	11.88

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	1,4-Benzenediol, 2,3-dichloro-5,6-d	312	C12H6Cl2N2O4	000000-00-0	1



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Vial: 5
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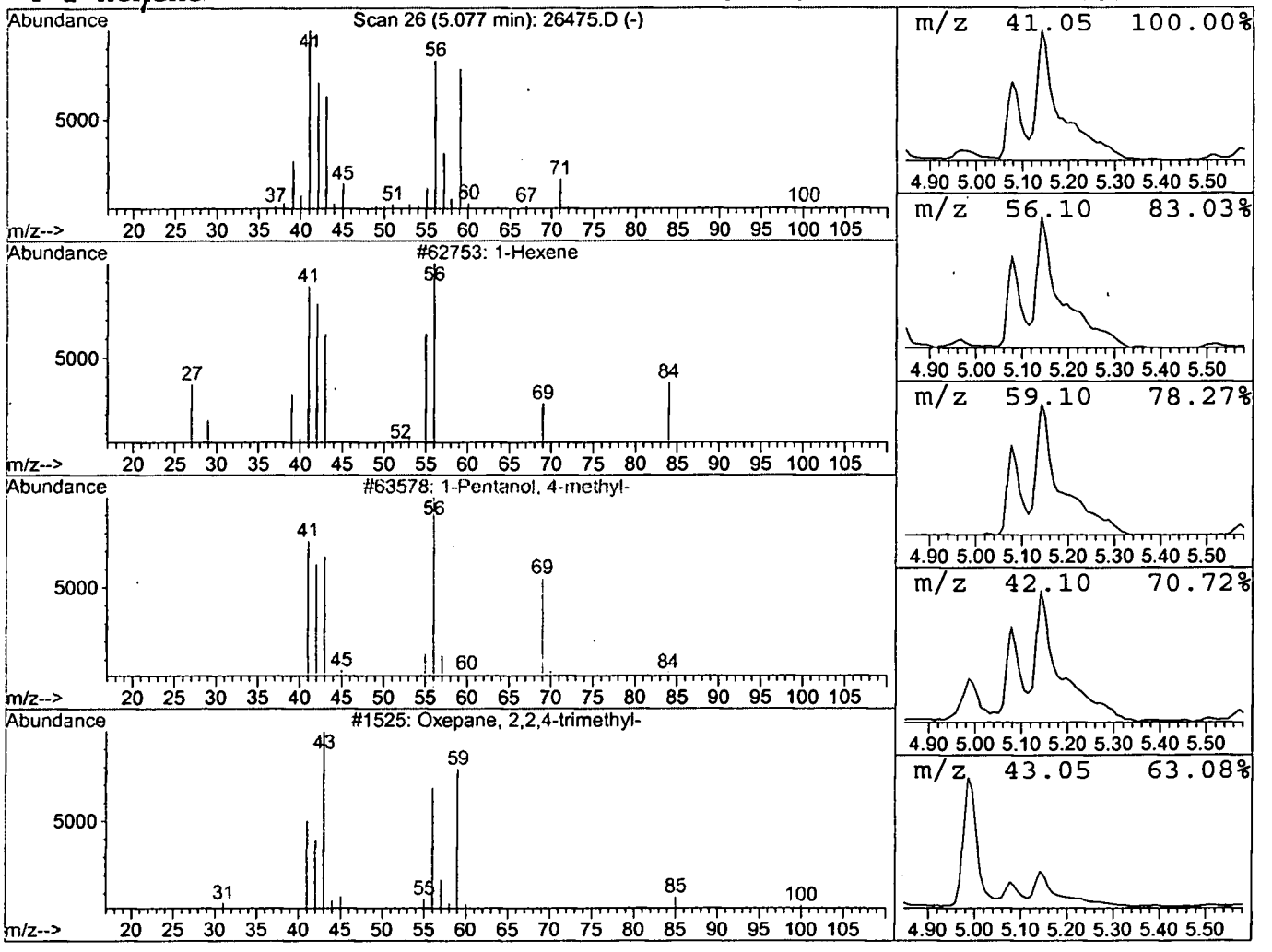
****

Peak Number 2 1-Hexene — Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.08	1.49 ng/uL	417216	1,4-Dichlorobenzene-d4	11.88

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Hexene	84	C6H12	000592-41-6	25
2	1-Pentanol, 4-methyl-	102	C6H14O	000626-89-1	25
3	Oxepane, 2,2,4-trimethyl-	100	C6H12O	023120-44-7	25
4	1-Hexene	84	C6H12	000592-41-6	25



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Acq On : 7 Dec 2001 9:45 pm
Sample : gc/ms unknown 11/6
Misc :
MS Integration Params: LSCINT.P

Vial: 5
Operator: GALOS
Inst : GC/MS Ins
Multiplr: 1.00

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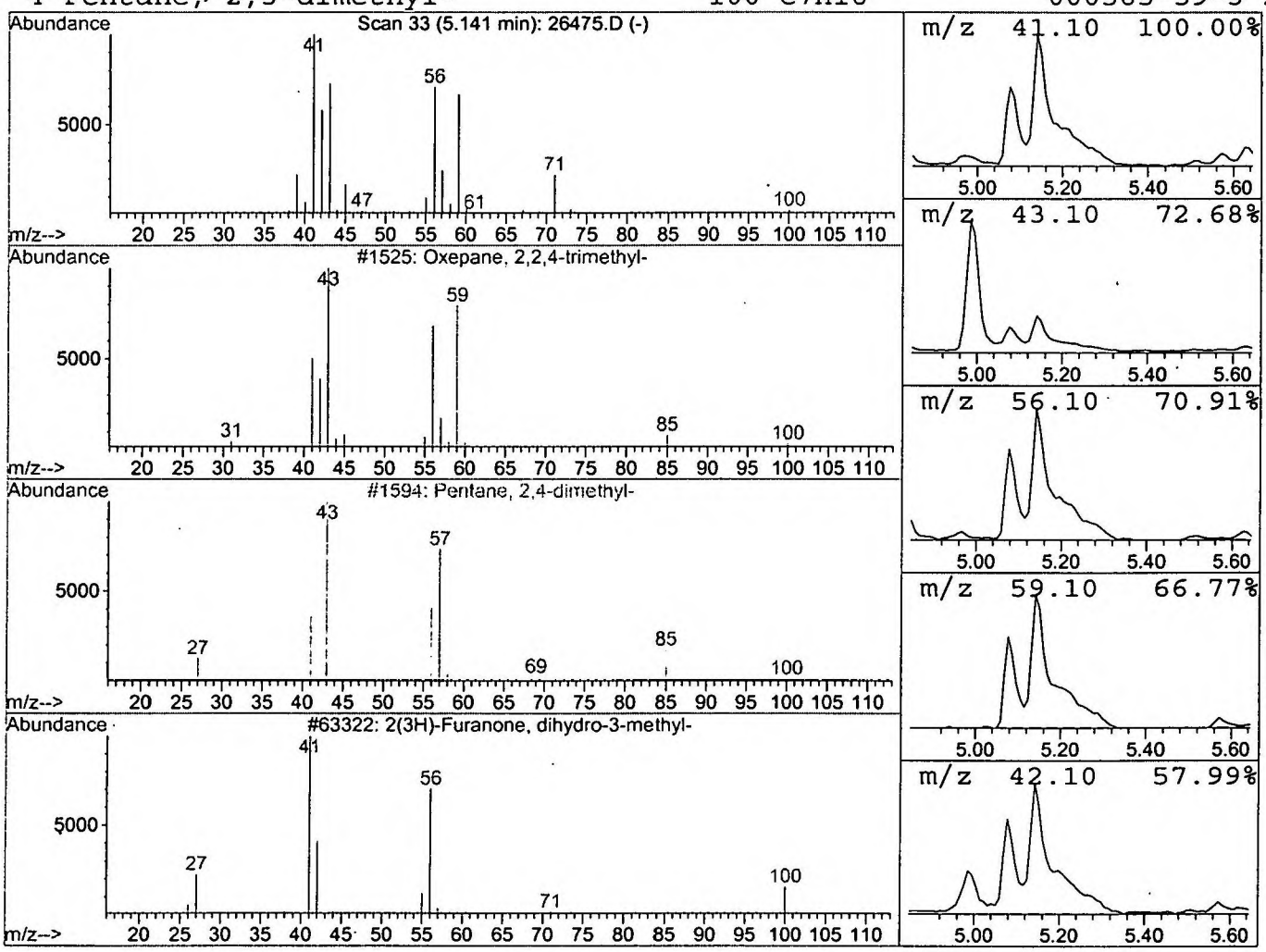
Peak Number 3 Oxepane, 2,2,4-trimethyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.14	4.43 ng/uL	1242160	1,4-Dichlorobenzene-d4	11.88

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Oxepane, 2,2,4-trimethyl-	100	C6H12O	023120-44-7	52
2	Pentane, 2,4-dimethyl-	100	C7H16	000108-08-7	32
3	2(3H)-Furanone, dihydro-3-methyl-	100	C5H8O2	001679-47-6	25
4	Pentane, 2,3-dimethyl-	100	C7H16	000565-59-3	25



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Misc :
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Vial: 5
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Multiplr: 1.00

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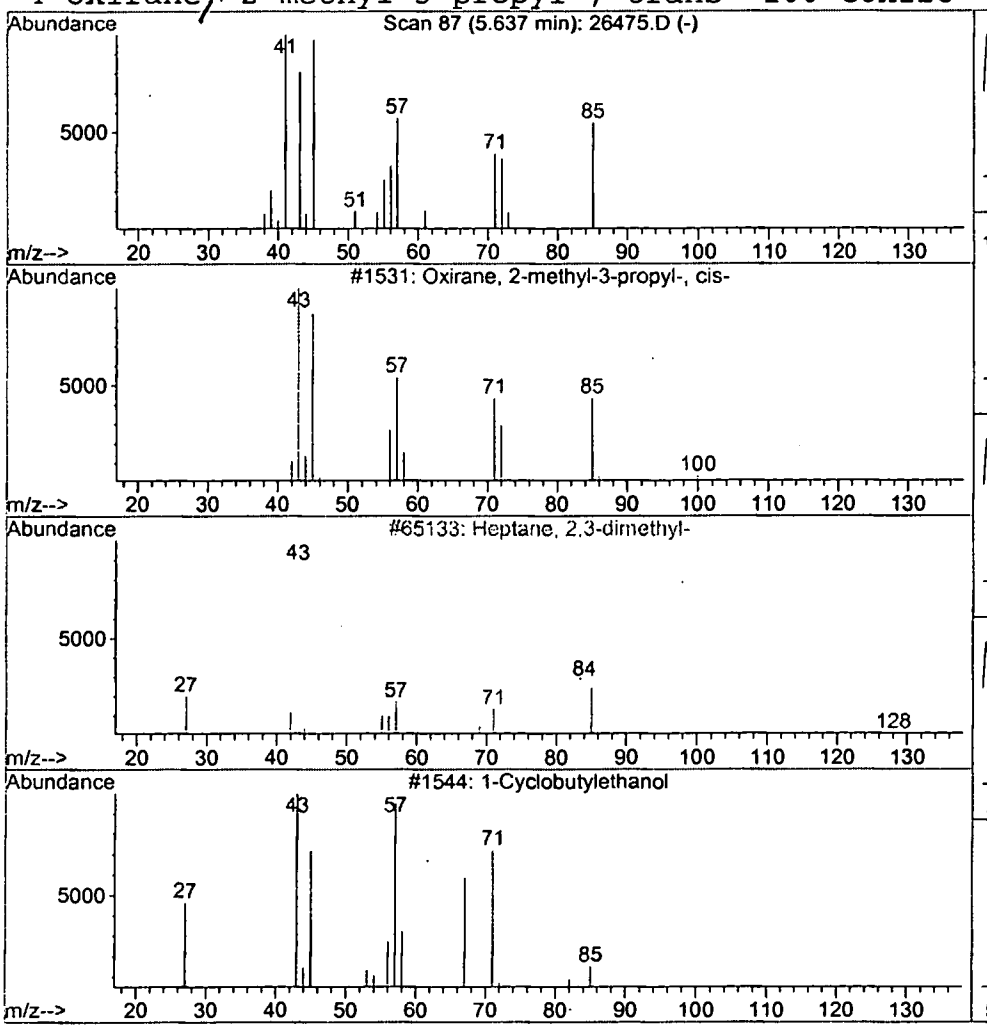
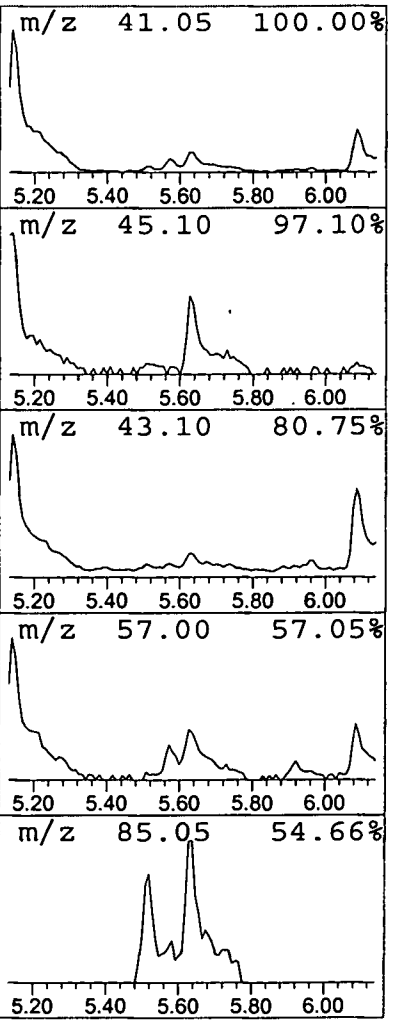
Peak Number 4 Oxirane, 2-methyl-3-propyl-, c Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.64	0.47 ng/uL	133131	1,4-Dichlorobenzene-d4	11.88

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Oxirane, 2-methyl-3-propyl-, cis-	100	C6H12O	006124-90-9	36
2	Heptane, 2,3-dimethyl-	128	C9H20	003074-71-3	17
3	1-Cyclobutylethanol	100	C6H12O	038401-41-1	17
4	Oxirane, 2-methyl-3-propyl-, trans-	100	C6H12O	006124-91-0	17





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Multiplr: 1.00

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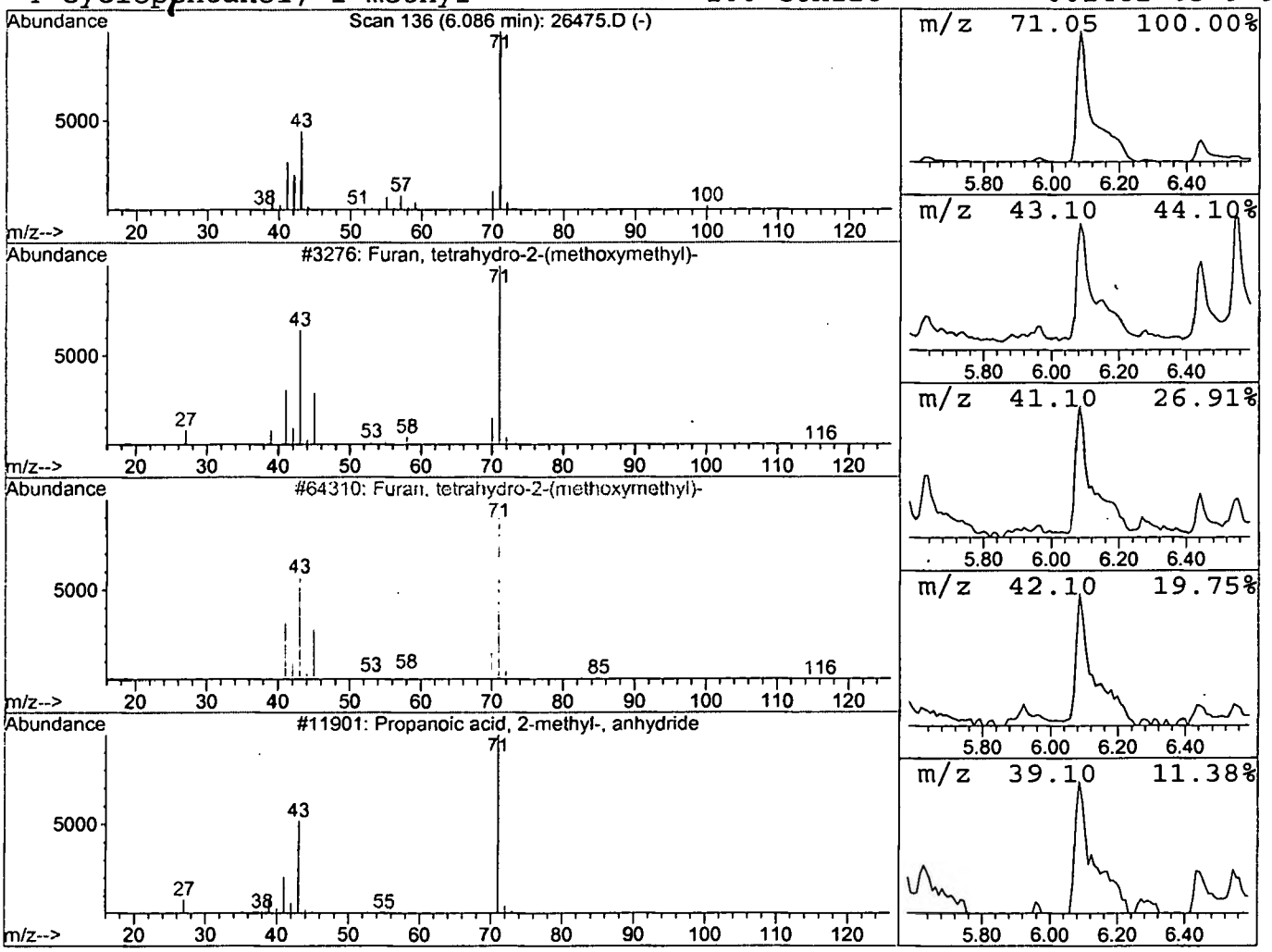
Peak Number 5 Furan, tetrahydro-2-(methoxyme Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.09	2.19 ng/uL	615331	1,4-Dichlorobenzene-d4	11.88

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Furan, tetrahydro-2-(methoxymethyl)	116	C6H12O2	019354-27-9	59
2	Furan, tetrahydro-2-(methoxymethyl)	116	C6H12O2	019354-27-9	45
3	Propanoic acid, 2-methyl-, anhydrid	158	C8H14O3	000097-72-3	28
4	Cyclopentanol, 1-methyl-	100	C6H12O	001462-03-9	9



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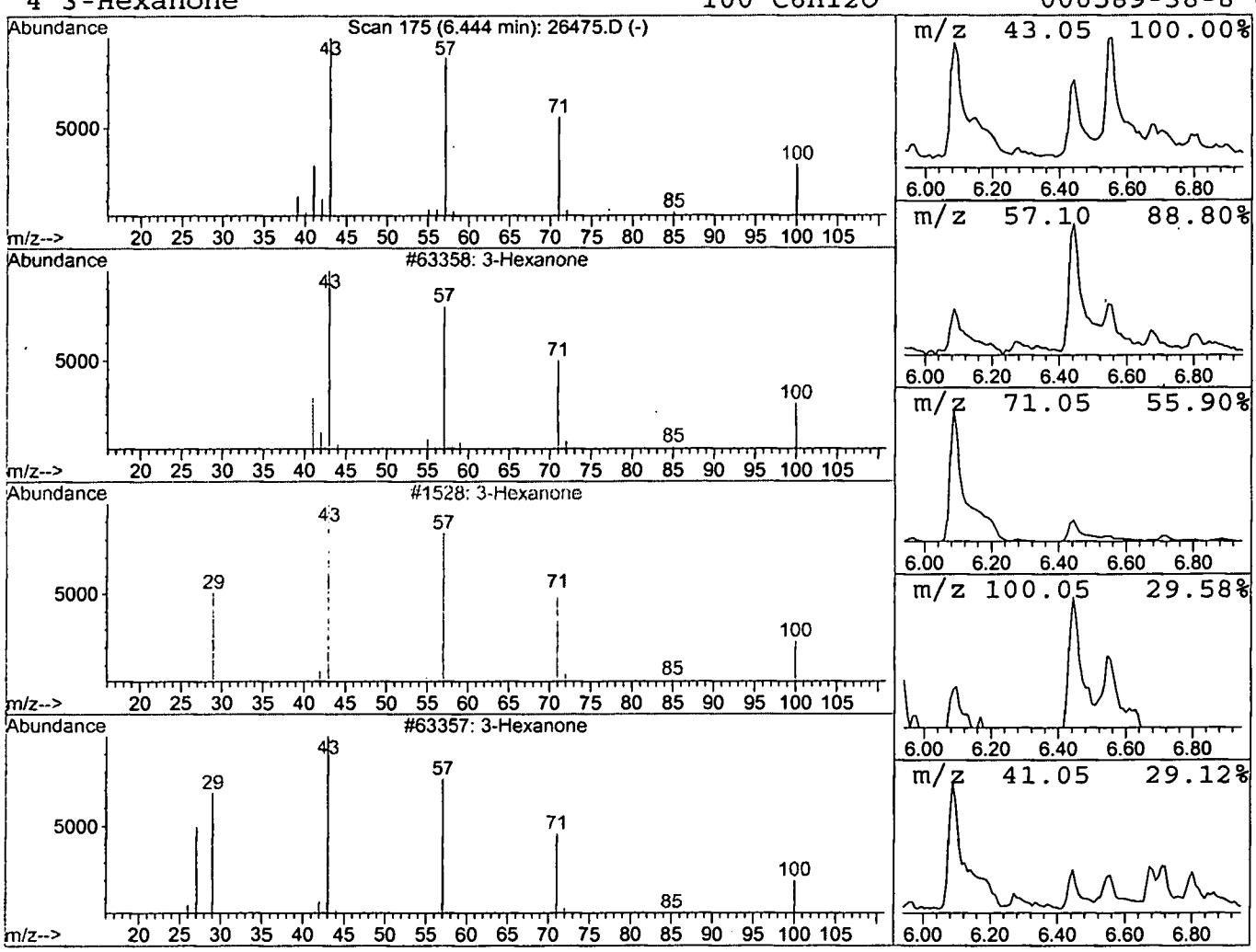
Library : C:\DATABASE\NBS75K.L

****

Peak Number 6 3-Hexanone — Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.44	0.67 ng/uL	187057	1,4-Dichlorobenzene-d4	11.88

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	90
4	3-Hexanone	100	C6H12O	000589-38-8	64



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Acq On : 7 Dec 2001 9:45 pm Operator: GALOS
Sample : gc/ms unknown 11/6 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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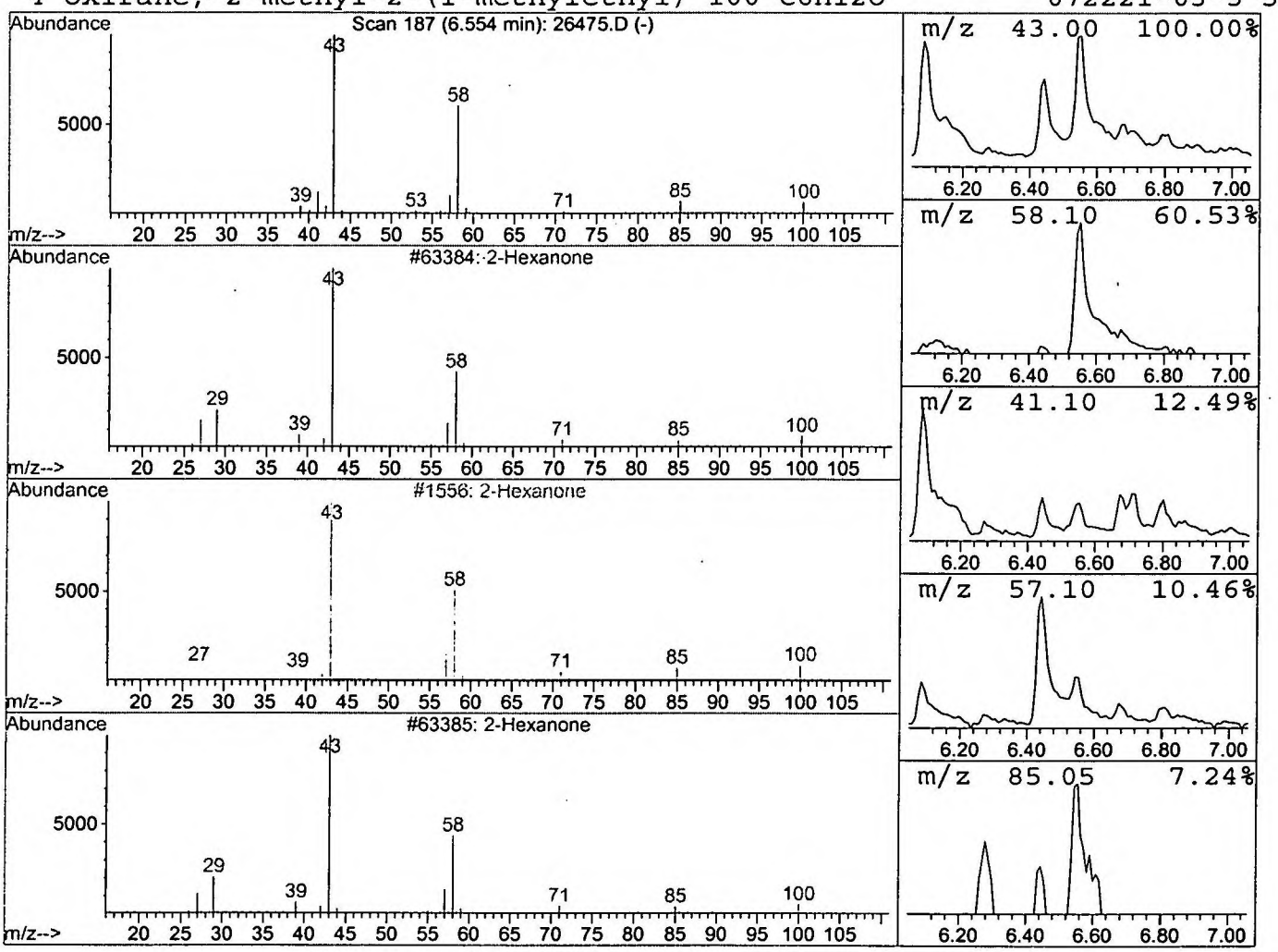
Peak Number 7 2-Hexanone Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.55	0.70 ng/uL	197213	1,4-Dichlorobenzene-d4	11.88

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Hexanone	100	C6H12O	000591-78-6	64
2	2-Hexanone	100	C6H12O	000591-78-6	53
3	2-Hexanone	100	C6H12O	000591-78-6	53
4	Oxirane, 2-methyl-2-(1-methylethyl)	100	C6H12O	072221-03-5	50



Library Search Compound Report

Data File : C:\HPCHEM\1\DATA\DEC0701\26475.D Vial: 5
Acq On : 7 Dec 2001 9:45 pm Operator: GALOS
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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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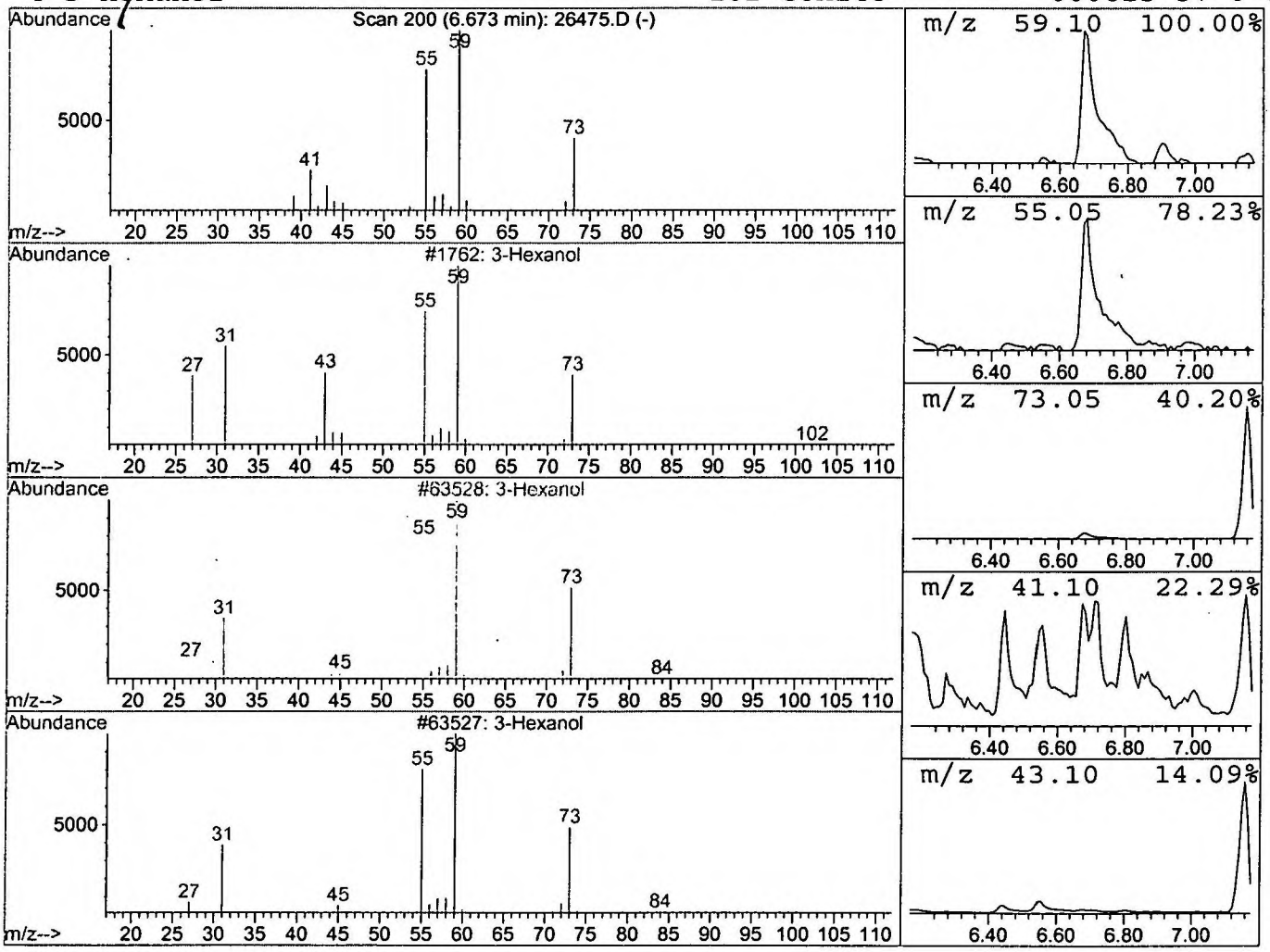
Peak Number 8 3-Hexanol Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.67	0.84 ng/uL	235131	1,4-Dichlorobenzene-d4	11.88

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	3-Hexanol	102	C6H14O	000623-37-0	78
2	3-Hexanol	102	C6H14O	000623-37-0	74
3	3-Hexanol	102	C6H14O	000623-37-0	74
4	3-Hexanol	102	C6H14O	000623-37-0	64



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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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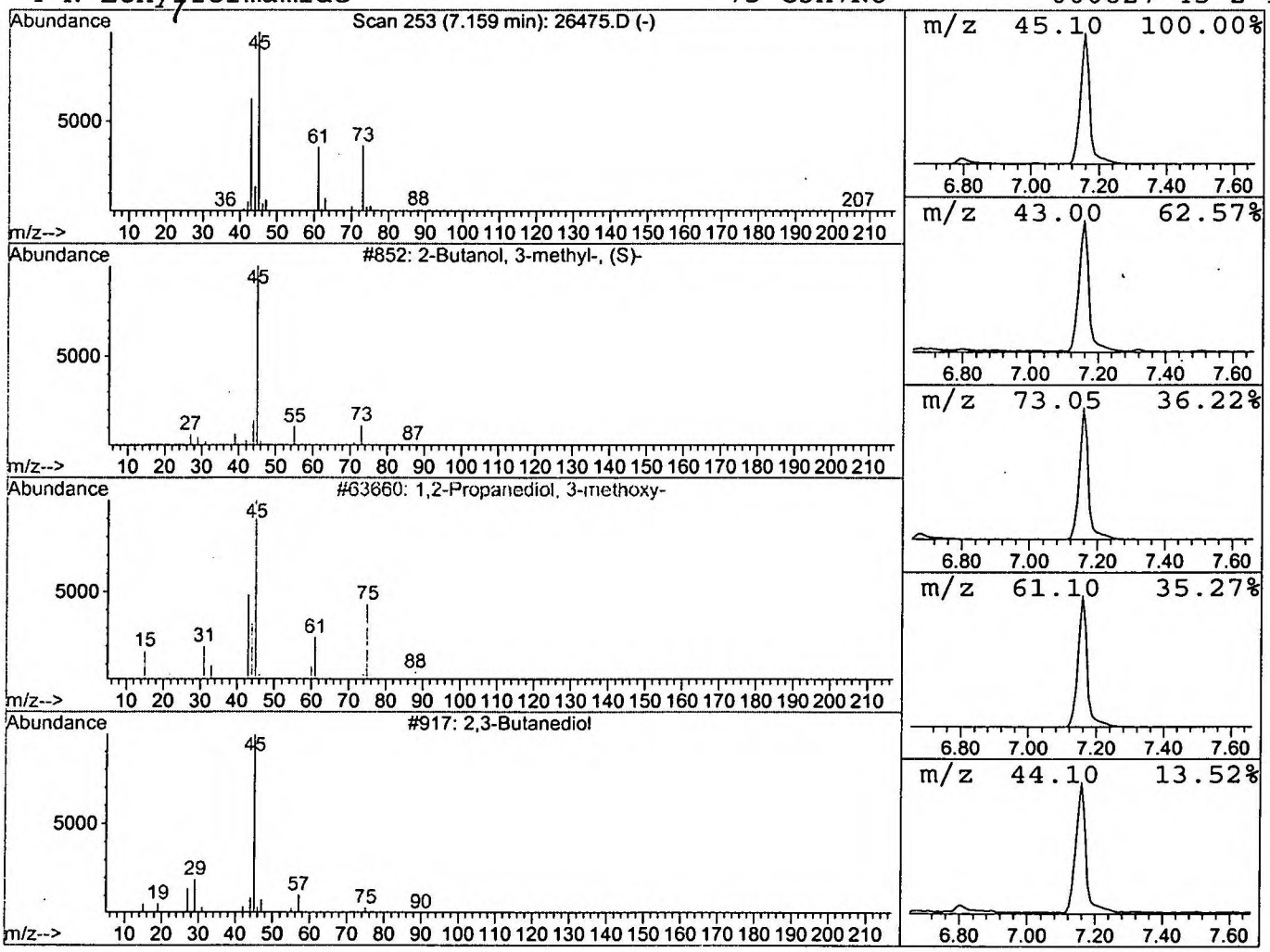
Peak Number 9 2-Butanol, 3-methyl-, (S)- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.16	12.21 ng/uL	3427830	1,4-Dichlorobenzene-d4	11.88

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Butanol, 3-methyl-, (S)-	88	C5H12O	001517-66-4	36
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



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Vial: 5
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

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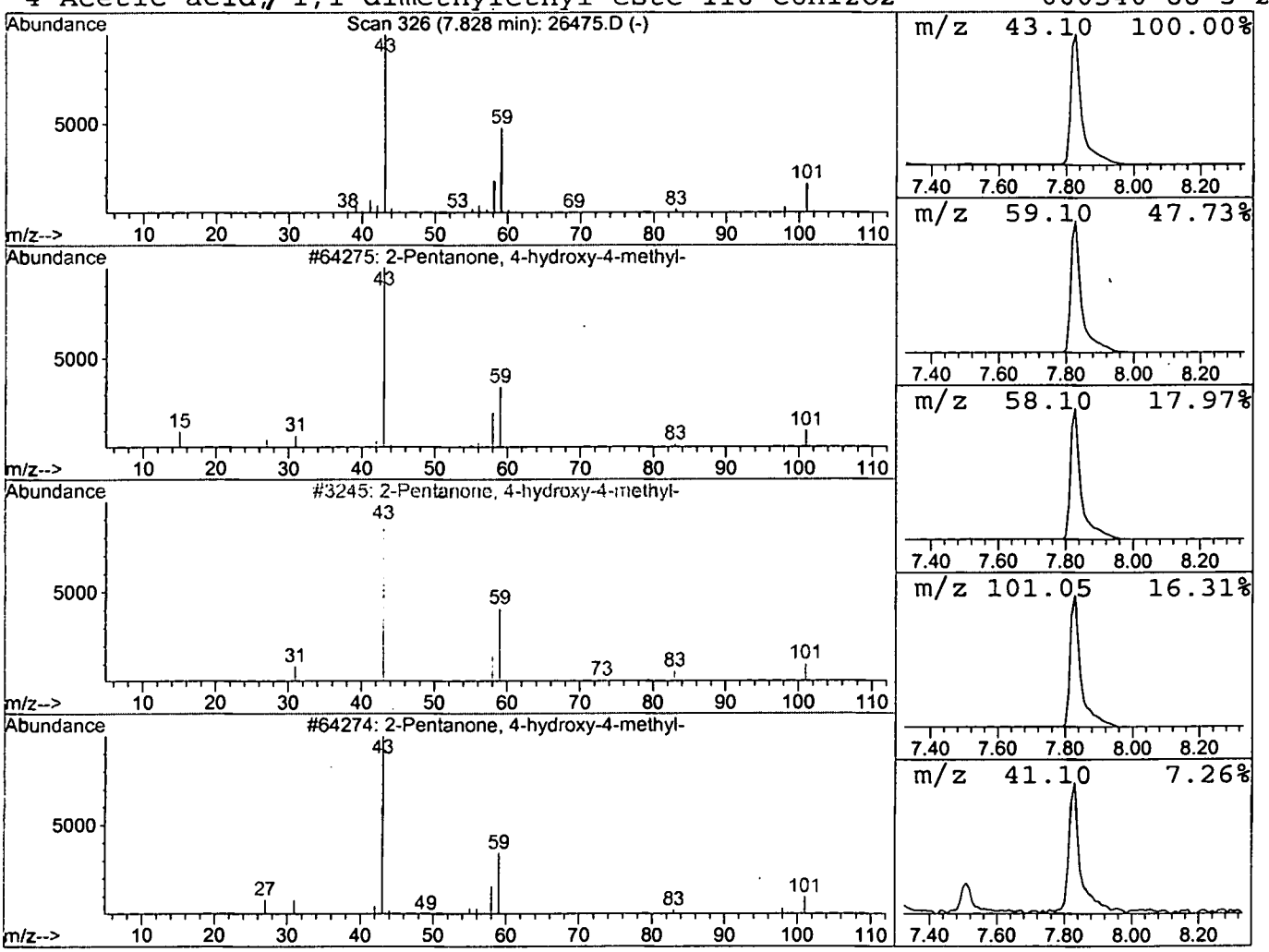
Peak Number 10 2-Pentanone, 4-hydroxy-4-methy Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.83	7.53 ng/uL	2113190	1,4-Dichlorobenzene-d4	11.88

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	40
4	Acetic acid, 1,1-dimethylethyl este	116	C6H12O2	000540-88-5	28



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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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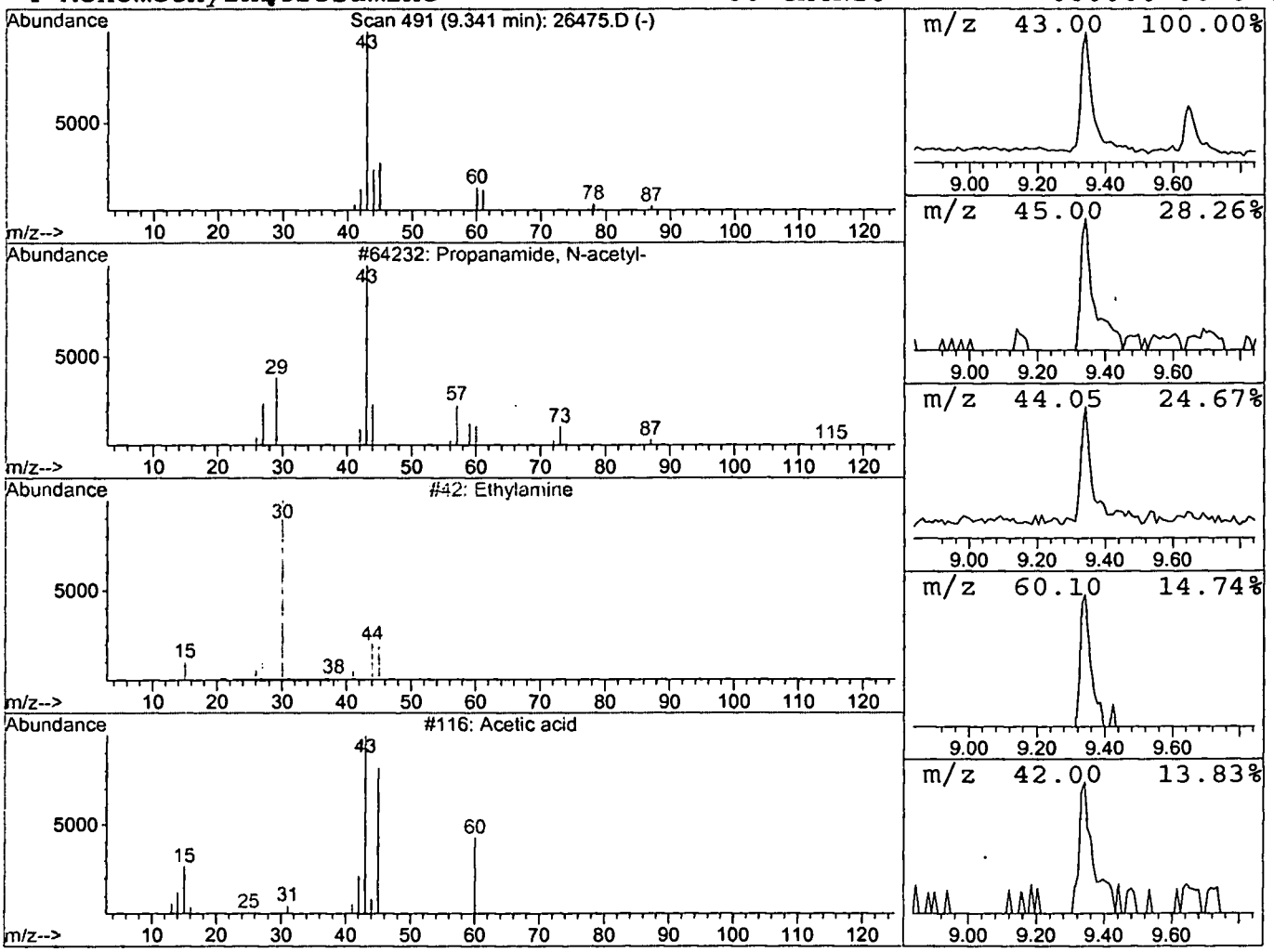
Peak Number 11 Propanamide, N-acetyl- Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.34	0.45 ng/uL	124989	1,4-Dichlorobenzene-d4	11.88

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Propanamide, N-acetyl-	115	C5H9NO2	019264-34-7	33
2	Ethylamine	45	C2H7N	000075-04-7	9
3	Acetic acid	60	C2H4O2	000064-19-7	7
4	Monomethylnitrosamine	60	CH4N2O	000000-00-0	7



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Misc :
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Vial: 5
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

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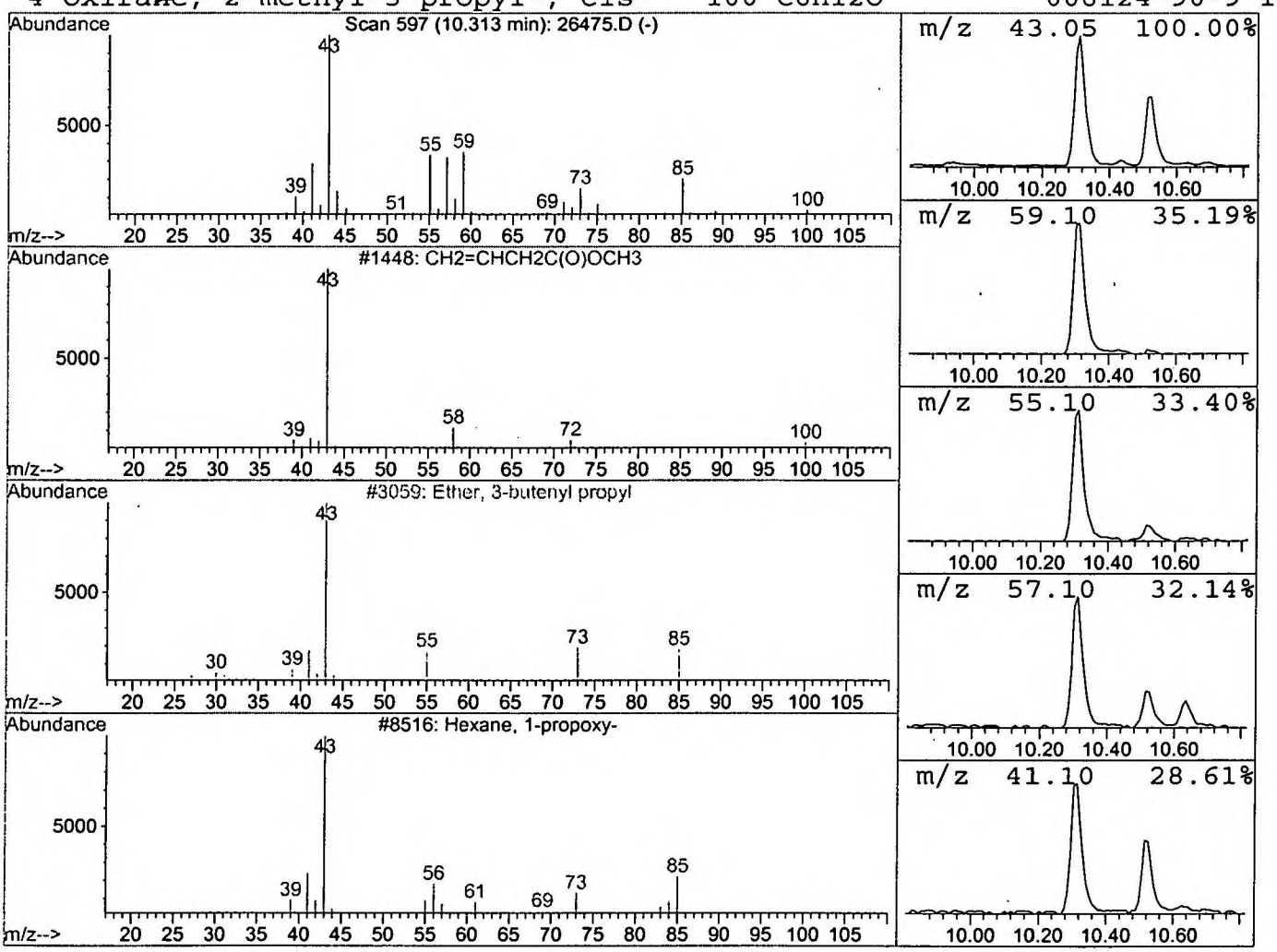
Peak Number 12 CH2=CHCH2C(O)OCH3 Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.31	2.93 ng/uL	822222	1,4-Dichlorobenzene-d4	11.88

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	CH2=CHCH2C(O)OCH3	100	C5H8O2	003724-55-8	27
2	Ether, 3-butenyl propyl	114	C7H14O	034061-75-1	23
3	Hexane, 1-propoxy-	144	C9H20O	053685-78-2	23
4	Oxirane, 2-methyl-3-propyl-, cis-	100	C6H12O	006124-90-9	17



Library Search Compound Report

Data File : C:\HPCHEM\1\DATA\DEC0701\26475.D
Acq On : 7 Dec 2001 9:45 pm
Sample : gc/ms unknown 11/6
Misc :
MS Integration Params: LSCINT.P

Vial: 5
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

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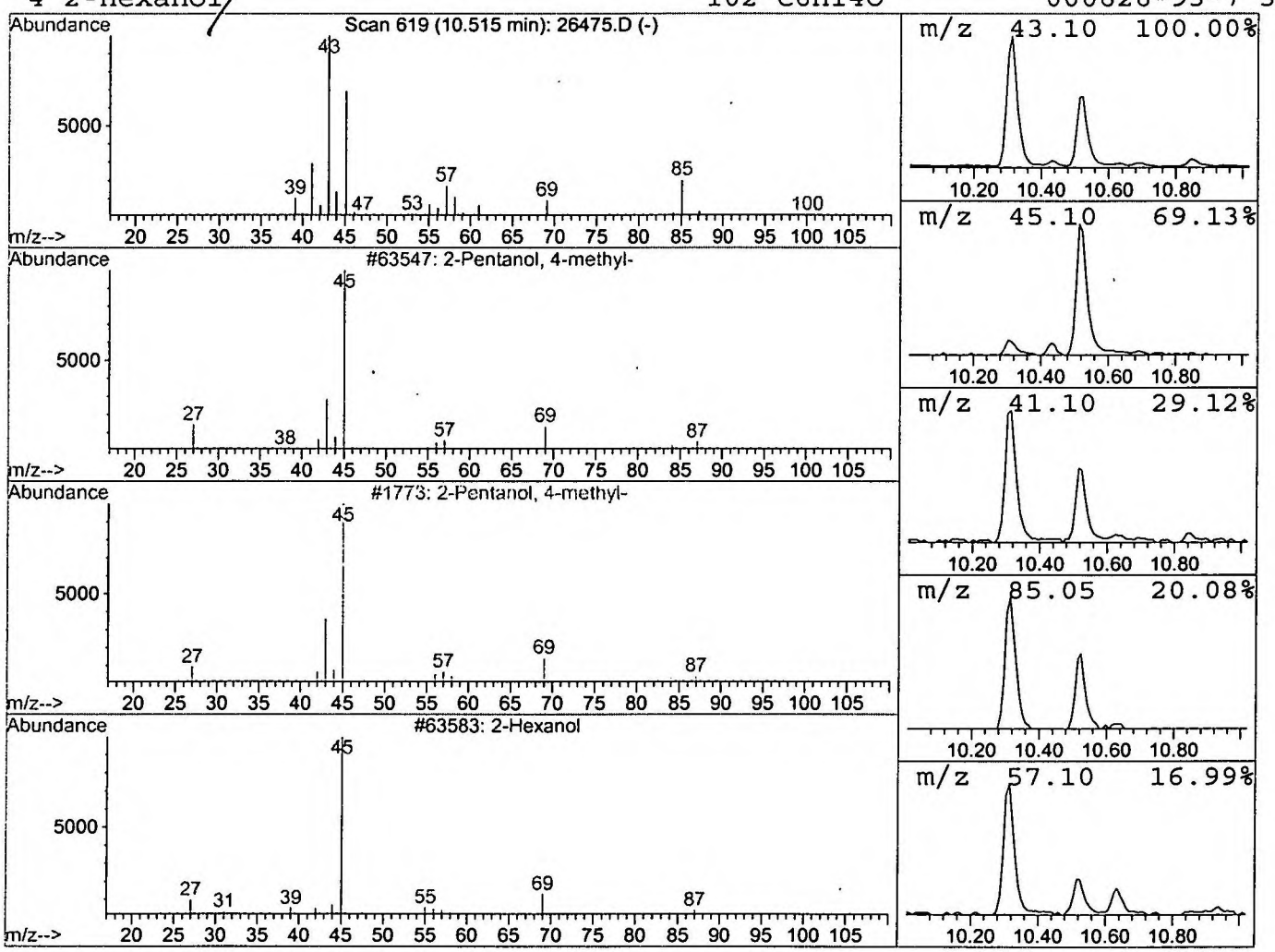
Peak Number 13 2-Pentanol, 4-methyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.52	1.41 ng/uL	396501	1,4-Dichlorobenzene-d4	11.88

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Pentanol, 4-methyl-	102	C6H14O	000108-11-2	38
2	2-Pentanol, 4-methyl-	102	C6H14O	000108-11-2	37
3	2-Hexanol	102	C6H14O	000626-93-7	35
4	2-Hexanol	102	C6H14O	000626-93-7	32



Tentatively Identified Compound (LSC) summary

Operator ID: GALOS Date Acquired: 7 Dec 2001 9:45 pm
Data File: C:\HPCHEM\1\DATA\DEC0701\26475.D
Name: gc/ms unknown 11/6
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	4.99	1.6	ng/uL	456682	ISTD01	11.88	5614080	20.0
1-Hexene	5.08	1.5	ng/uL	417216	ISTD01	11.88	5614080	20.0
Oxepane, 2,2,4-trime	5.14	4.4	ng/uL	1242160	ISTD01	11.88	5614080	20.0
Oxirane, 2-methyl-3-	5.64	0.5	ng/uL	133131	ISTD01	11.88	5614080	20.0
Furan, tetrahydro-2-	6.09	2.2	ng/uL	615331	ISTD01	11.88	5614080	20.0
3-Hexanone	6.44	0.7	ng/uL	187057	ISTD01	11.88	5614080	20.0
2-Hexanone	6.55	0.7	ng/uL	197213	ISTD01	11.88	5614080	20.0
3-Hexanol	6.67	0.8	ng/uL	235131	ISTD01	11.88	5614080	20.0
2-Butanol, 3-methyl-	7.16	12.2	ng/uL	3427830	ISTD01	11.88	5614080	20.0
2-Pentanone, 4-hydro	7.83	7.5	ng/uL	2113190	ISTD01	11.88	5614080	20.0
Propanamide, N-acety	9.34	0.4	ng/uL	124989	ISTD01	11.88	5614080	20.0
CH2=CHCH2C(O)OCH3	10.31	2.9	ng/uL	822222	ISTD01	11.88	5614080	20.0
2-Pentanol, 4-methyl	10.52	1.4	ng/uL	396501	ISTD01	11.88	5614080	20.0

26475.D NP112701.M Thu Dec 13 11:19:48 2001