User: Galos, Marie
Westchester Dept. of Labs & Research
Date: 12/04/2001 Time: 15:11:16

Sample: AD26388
Analysis: $GCMS GCMS Scan For Unknowns
Units: ug
Started: 12/4/01 14:38 Ended: 12/4/01 14:38 Analyst: MG
Page: 1

	Component Name	Viol	Result
1	Other unknown compounds		see comment

Comments for Sample AD26388 - Analysis $GCMS

Westchester Dept. of Labs & Research
User: Galos, Marie
Date: 12-04-2001 Time: 15:11:27

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.

(QT Reviewed)

Quantitation Report (QT Reviewed)

Data File : C:\HPCHEM\1\DATA\NOV3001\26388.D Vial: 14
Acq On : 1 Dec 2001 12:02 am 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	1448612	20.00	ng/uL	-0.05
19) Naphthalene-d8	15.50	136	5622454	20.00	ng/uL	-0.05
31) Acenaphthene-d10	20.78	164	2508790	20.00	ng/uL	-0.04
49) Phenanthrene-d10	25.19	188	3641229	20.00	ng/uL	-0.05
59) Chrysene-d12	33.19	240	2317715	20.00	ng/uL	-0.06
68) Perylene-d12	37.28	264	1660070	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	11.15	99	884	0.01	ng/uL	0.02
Spiked Amount 75.000	Range 19 - 32		Recovery =	0.01%#		
6) 2-Chlorophenol-d4	11.90	132	1567	0.02	ng/uL	0.47
Spiked Amount 75.000	Range 34 - 84		Recovery =	0.03%#		
7) 1,2-Dichlorobenzene-d4	12.41	152	14251	0.23	ng/uL	-0.05
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	4818	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	12.90	45	3149	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
26388.D 12701.M Tue Dec 04 08:19:40 2001 Page 1

Quantitation Report

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Acq On : 1 Dec 2001 12:02 am Operator: GALOS
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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

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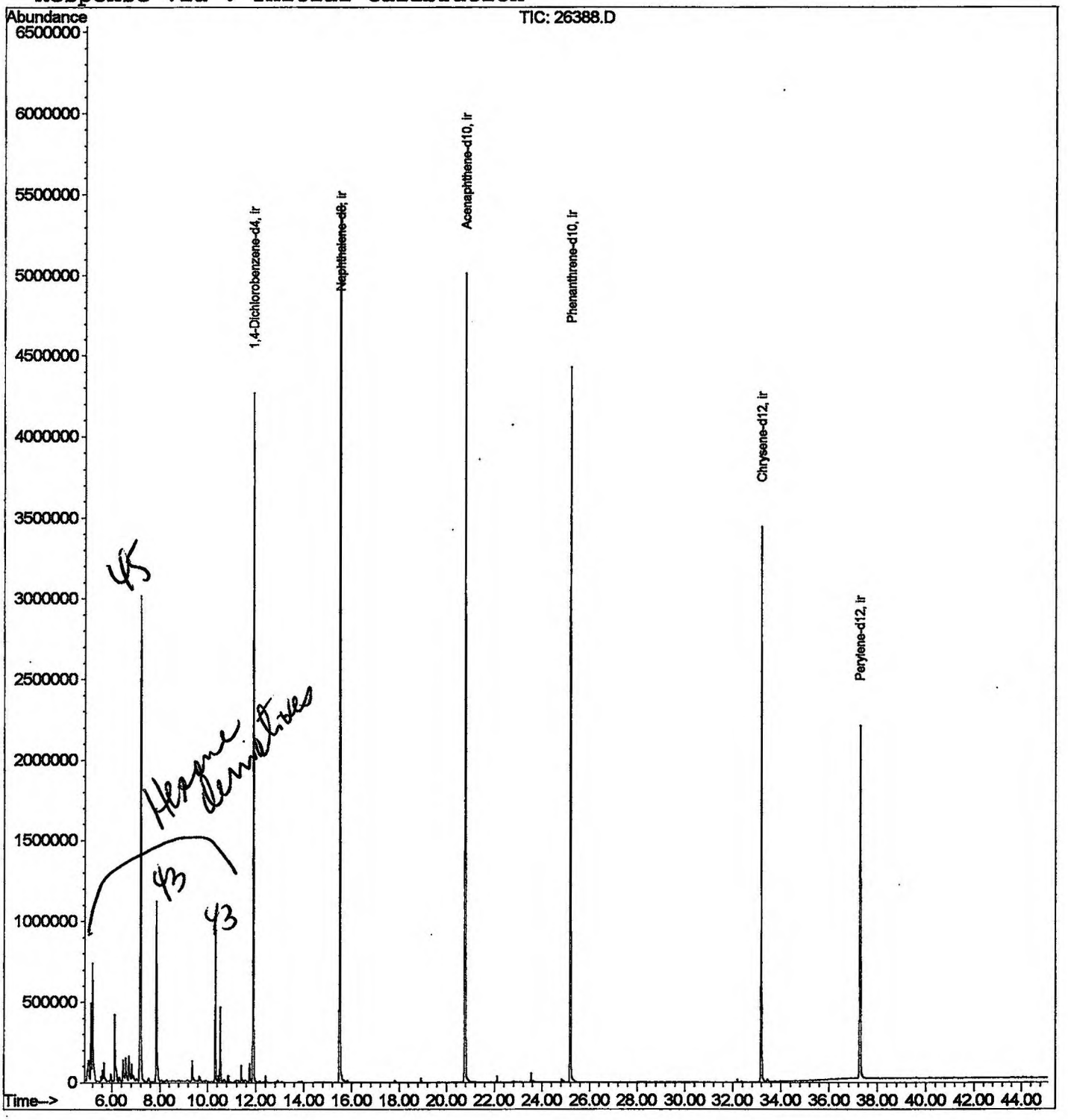


LSC Area Percent Report

Data File : C:\HPCHEM\1\DATA\NOV3001\26388.D Vial: 14
Acq On : 1 Dec 2001 12:02 am 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

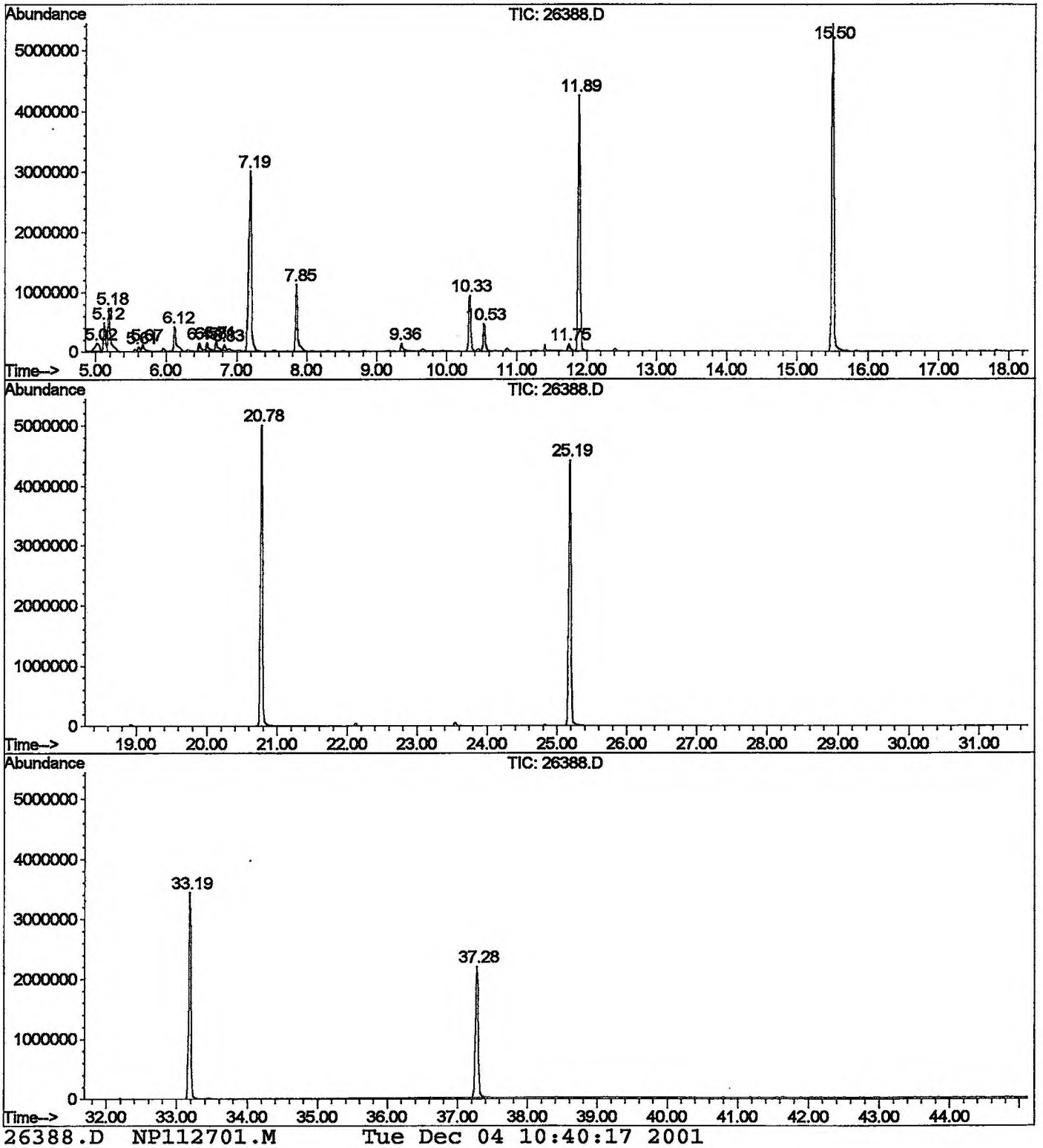
peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	peak area	peak % max.	% of total
1	5.020	9	19	27	rBV	125680	591056	5.24%	0.801%
2	5.121	27	30	34	rVV	484192	865260	7.67%	1.172%
3	5.185	34	37	55	rVB	732758	1753565	15.55%	2.375%
4	5.607	80	83	87	rVV	68585	141844	1.26%	0.192%
5	5.671	87	90	103	rVB	113592	270347	2.40%	0.366%
6	6.120	135	139	156	rBV	413000	1027333	9.11%	1.391%
7	6.469	173	177	182	rBV	129045	256746	2.28%	0.348%
8	6.579	186	189	199	rVB	136598	304613	2.70%	0.413%
9	6.707	199	203	212	rBV	150300	370594	3.29%	0.502%
10	6.826	212	216	220	rBV	91688	175190	1.55%	0.237%
11	7.193	248	256	274	rVB	3008000	8078196	71.64%	10.941%
12	7.853	324	328	345	rBV	1117721	2350041	20.84%	3.183%
13	9.357	488	492	505	rBV	125624	281259	2.49%	0.381%
14	10.329	593	598	607	rBV	940142	1975687	17.52%	2.676%
15	10.531	615	620	631	rBV	464985	947159	8.40%	1.283%
16	11.750	749	753	761	rBV2	114263	251345	2.23%	0.340%
17	11.888	763	768	782	rVB	4262671	8792768	77.98%	11.909%
18	15.501	1156	1162	1181	rBV	5442676	11275706	100.00%	15.272%
19	20.783	1731	1738	1756	rBV	5011837	10849790	96.22%	14.695%
20	25.194	2212	2219	2230	rBV2	4424170	9900458	87.80%	13.409%
21	33.190	3084	3091	3107	rBV2	3439298	7434469	65.93%	10.069%
22	37.280	3529	3537	3548	rBV2	2183682	5938537	52.67%	8.043%

Sum of corrected areas: 73831963

26388.D NP112701.M Tue Dec 04 10:40:16 2001

LSC Report - Integrated Chromatogram

File : C:\HPCHEM\1\DATA\NOV3001\26388.D
Operator : GALOS
Acquired : 1 Dec 2001 12:02 am using AcqMethod NP112701
Instrument : GC/MS Ins
Sample Name: gc/ms unknown 11/3
Misc Info : .
Vial Number: 14
Quant File :NP112701.RES (RTE Integrator)



26388.D NP112701.M Tue Dec 04 10:40:17 2001

Library Search Compound Report

Data File : C:\HPCHEM\1\DATA\NOV3001\26388.D
Acq On : 1 Dec 2001 12:02 am
Sample : gc/ms unknown 11/3
Misc :
MS Integration Params: LSCINT.P

Vial: 14
Operator: GALOS
Inst : GC/MS Ins
Multiplr: 1.00

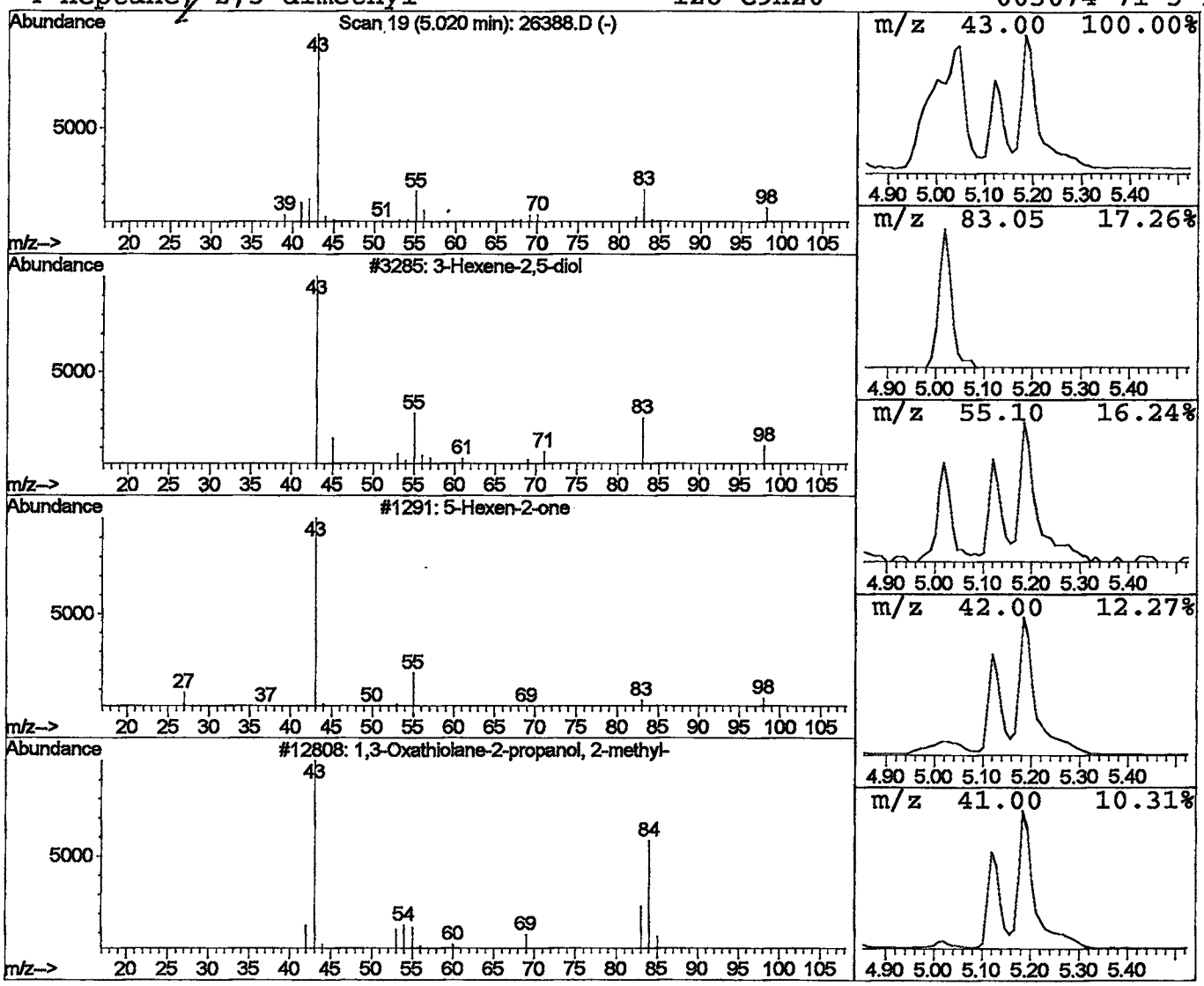
Quant Method : C:\HPCHEM\1\METHODS\NP112701.M (RTE Integrator)
Title : 208 625/TCLP calibration table
Library : C:\DATABASE\NBS75K.L

****

Peak Number 1 3-Hexene-2,5-diol Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.02	1.34 ng/uL	591056	1,4-Dichlorobenzene-d4	11.89

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	3-Hexene-2,5-diol	116	C6H12O2	007319-23-5	25
2	5-Hexen-2-one	98	C6H10O	000109-49-9	16
3	1,3-Oxathiolane-2-propanol, 2-methy	162	C7H14O2S	036651-30-6	9
4	Heptane, 2,3-dimethyl-	128	C9H20	003074-71-3	9



Library Search Compound Report

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

Vial: 14
Operator: GALOS
Inst : GC/MS Ins
Multiplr: 1.00

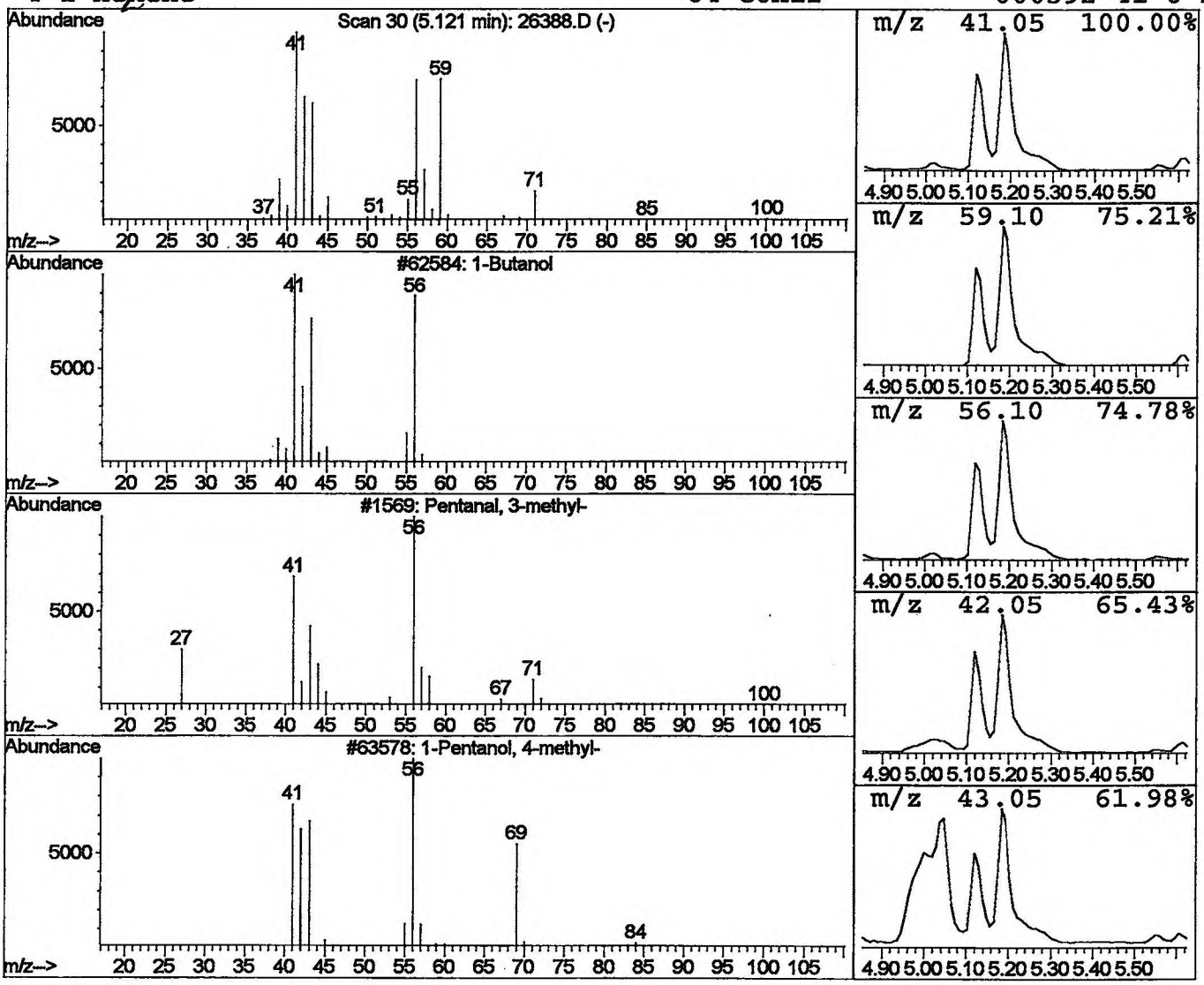
Quant Method : C:\HPCHEM\1\METHODS\NP112701.M (RTE Integrator)
Title : 208 625/TCLP calibration table
Library : C:\DATABASE\NBS75K.L

****

Peak Number 2 1-Butanol — Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.12	1.97 ng/uL	865260	1,4-Dichlorobenzene-d4	11.89

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Butanol	74	C4H10O	000071-36-3	37
2	Pentanal, 3-methyl-	100	C6H12O	015877-57-3	35
3	1-Pentanol, 4-methyl-	102	C6H14O	000626-89-1	28
4	1-Hexene	84	C6H12	000592-41-6	25



Library Search Compound Report

Data File : C:\HPCHEM\1\DATA\NOV3001\26388.D Vial: 14
Acq On : 1 Dec 2001 12:02 am Operator: GALOS
Sample : gc/ms unknown 11/3 Inst : GC/MS Ins
Misc : Multiplr: 1.00
MS Integration Params: LSCINT.P

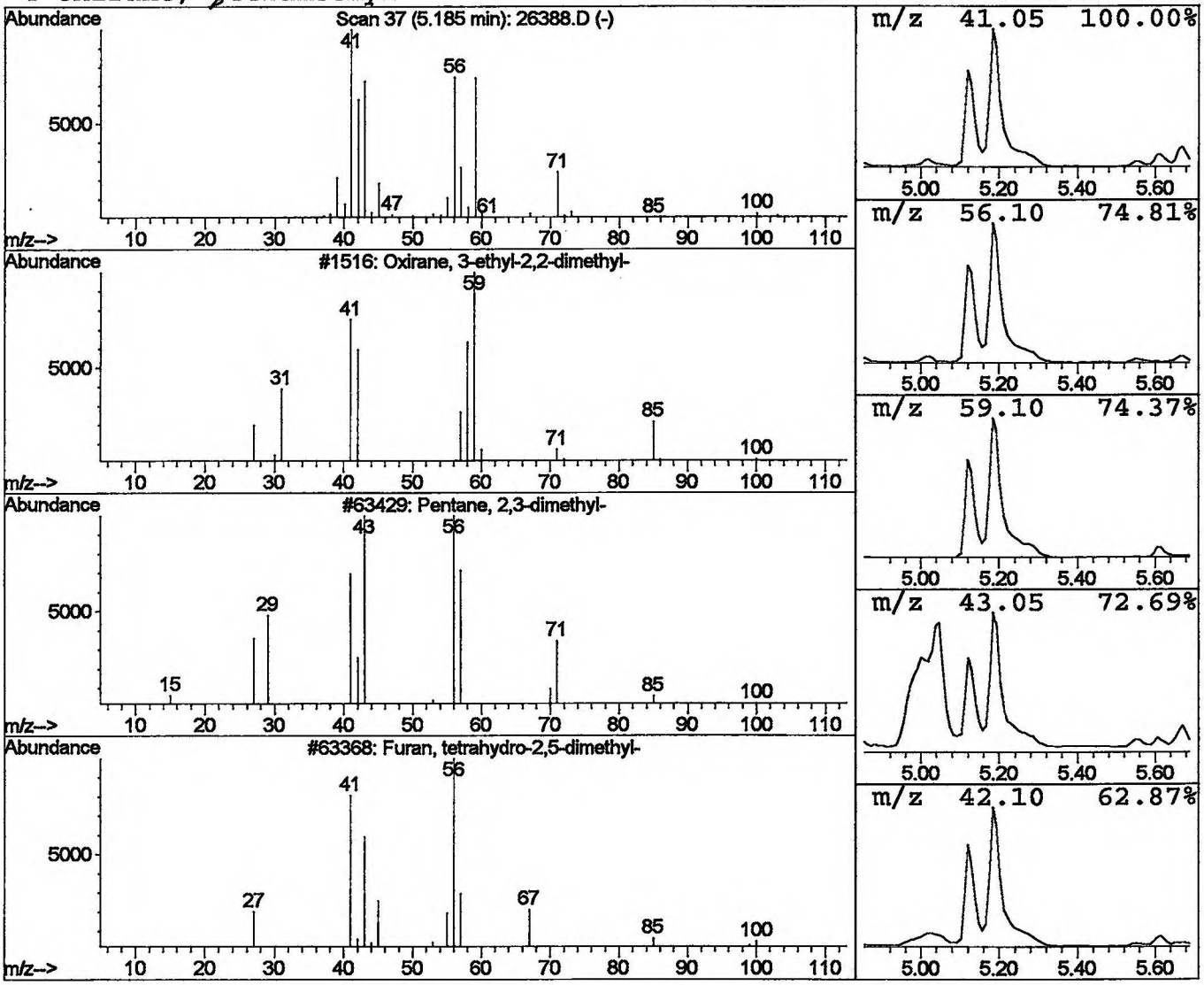
Quant Method : C:\HPCHEM\1\METHODS\NP112701.M (RTE Integrator)
Title : 208 625/TCLP calibration table
Library : C:\DATABASE\NBS75K.L

****

Peak Number 3 Oxirane, 3-ethyl-2,2-dimethyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.18	3.99 ng/uL	1753570	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	38
2	Pentane, 2,3-dimethyl-	100	C7H16	000565-59-3	35
3	Furan, tetrahydro-2,5-dimethyl-	100	C6H12O	001003-38-9	27
4	Oxirane, tetramethyl-	100	C6H12O	005076-20-0	27



Library Search Compound Report

Data File : C:\HPCHEM\1\DATA\NOV3001\26388.D
Acq On : 1 Dec 2001 12:02 am
Sample : gc/ms unknown 11/3
Misc :
MS Integration Params: LSCINT.P

Vial: 14
Operator: GALOS
Inst : GC/MS Ins
Multiplr: 1.00

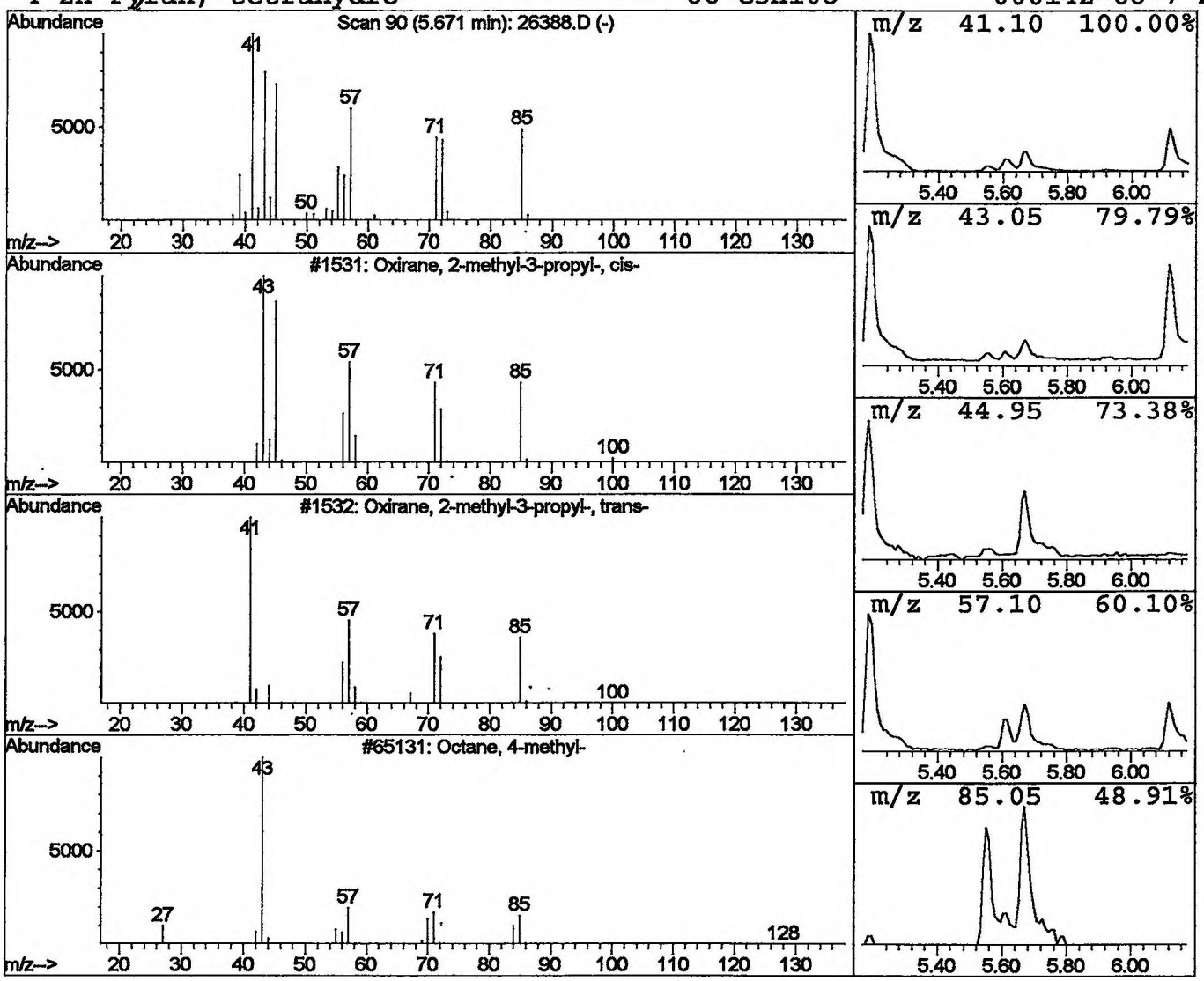
Quant Method : C:\HPCHEM\1\METHODS\NP112701.M (RTE Integrator)
Title : 208 625/TCLP calibration table
Library : C:\DATABASE\NBS75K.L

****

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.67	0.61 ng/uL	270347	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	42
2	Oxirane, 2-methyl-3-propyl-, trans-	100	C6H12O	006124-91-0	37
3	Octane, 4-methyl-	128	C9H20	002216-34-4	32
4	2H-Pyran, tetrahydro-	86	C5H10O	000142-68-7	23



Library Search Compound Report

Data File : C:\HPCHEM\1\DATA\NOV3001\26388.D
Acq On : 1 Dec 2001 12:02 am
Sample : gc/ms unknown 11/3
Misc :
MS Integration Params: LSCINT.P

Vial: 14
Operator: GALOS
Inst : GC/MS Ins
Multiplr: 1.00

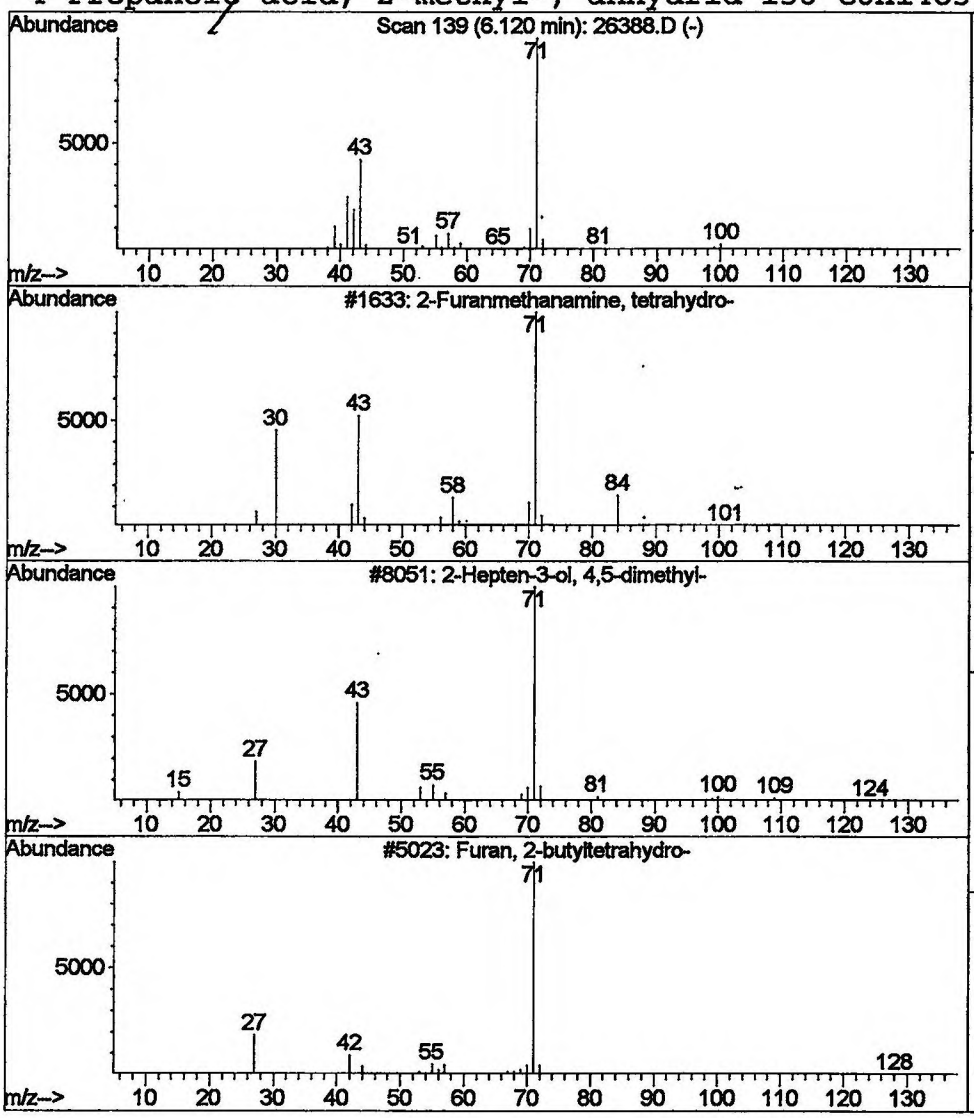
Quant Method : C:\HPCHEM\1\METHODS\NP112701.M (RTE Integrator)
Title : 208 625/TCLP calibration table
Library : C:\DATABASE\NBS75K.L

****

Peak Number 5 2-Furanmethanamine, tetrahydro Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.12	2.34 ng/uL	1027330	1,4-Dichlorobenzene-d4	11.89

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Furanmethanamine, tetrahydro-	101	C5H11NO	004795-29-3	40
2	2-Hepten-3-ol, 4,5-dimethyl-	142	C9H18O	055956-37-1	40
3	Furan, 2-butyltetrahydro-	128	C8H16O	001004-29-1	39
4	Propanoic acid, 2-methyl-, anhydrid	158	C8H14O3	000097-72-3	38



Library Search Compound Report

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

Vial: 14
Operator: GALOS
Inst : GC/MS Ins
Multiplr: 1.00

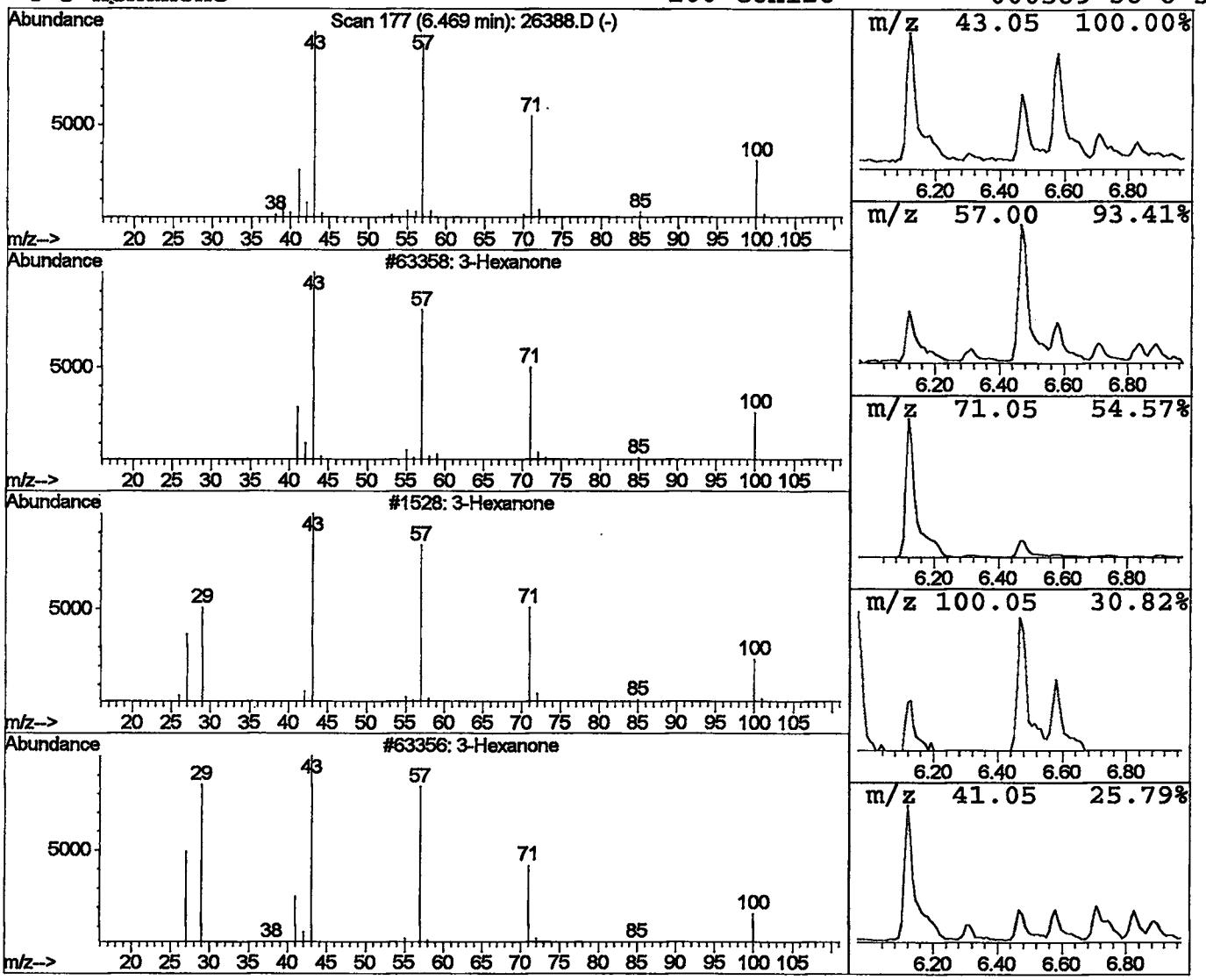
Quant Method : C:\HPCHEM\1\METHODS\NP112701.M (RTE Integrator)
Title : 208 625/TCLP calibration table
Library : C:\DATABASE\NBS75K.L

****

Peak Number 6 3-Hexanone Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.47	0.58 ng/uL	256746	1,4-Dichlorobenzene-d4	11.89

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	3-Hexanone	100	C6H12O	000589-38-8	87
2	3-Hexanone	100	C6H12O	000589-38-8	86
3	3-Hexanone	100	C6H12O	000589-38-8	78
4	3-Hexanone	100	C6H12O	000589-38-8	56



Library Search Compound Report

Data File : C:\HPCHEM\1\DATA\NOV3001\26388.D
Acq On : 1 Dec 2001 12:02 am
Sample : gc/ms unknown 11/3
Misc :
MS Integration Params: LSCINT.P

Vial: 14
Operator: GALOS
Inst : GC/MS Ins
Multiplr: 1.00

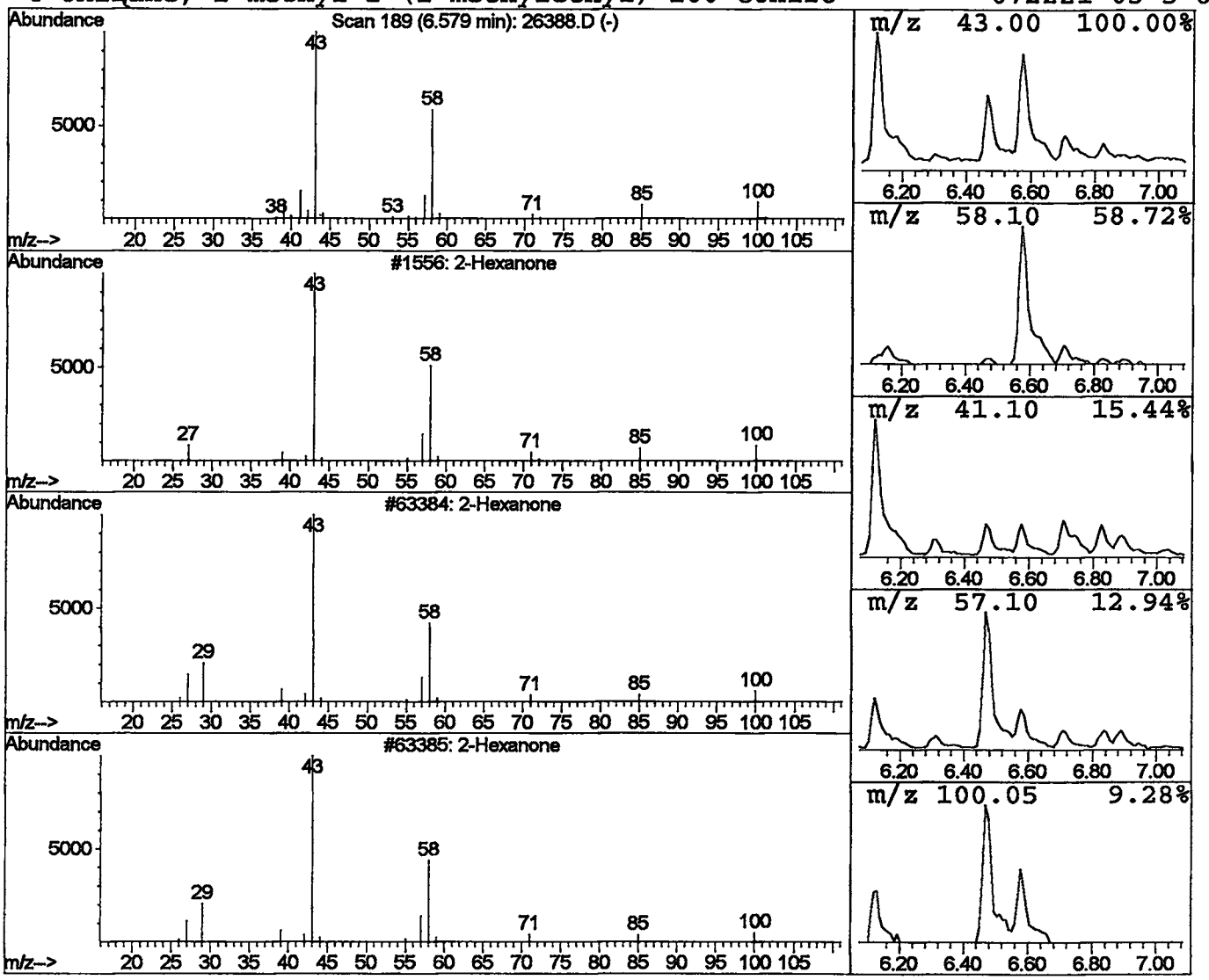
Quant Method : C:\HPCHEM\1\METHODS\NP112701.M (RTE Integrator)
Title : 208 625/TCLP calibration table
Library : C:\DATABASE\NBS75K.L

****

Peak Number 7 2-Hexanone — Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.58	0.69 ng/uL	304613	1,4-Dichlorobenzene-d4	11.89

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Hexanone	100	C6H12O	000591-78-6	90
2	2-Hexanone	100	C6H12O	000591-78-6	72
3	2-Hexanone	100	C6H12O	000591-78-6	72
4	Oxirane, 2-methyl-2-(1-methylethyl)	100	C6H12O	072221-03-5	64



Library Search Compound Report

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

Vial: 14
Operator: GALOS
Inst : GC/MS Ins
Multiplr: 1.00

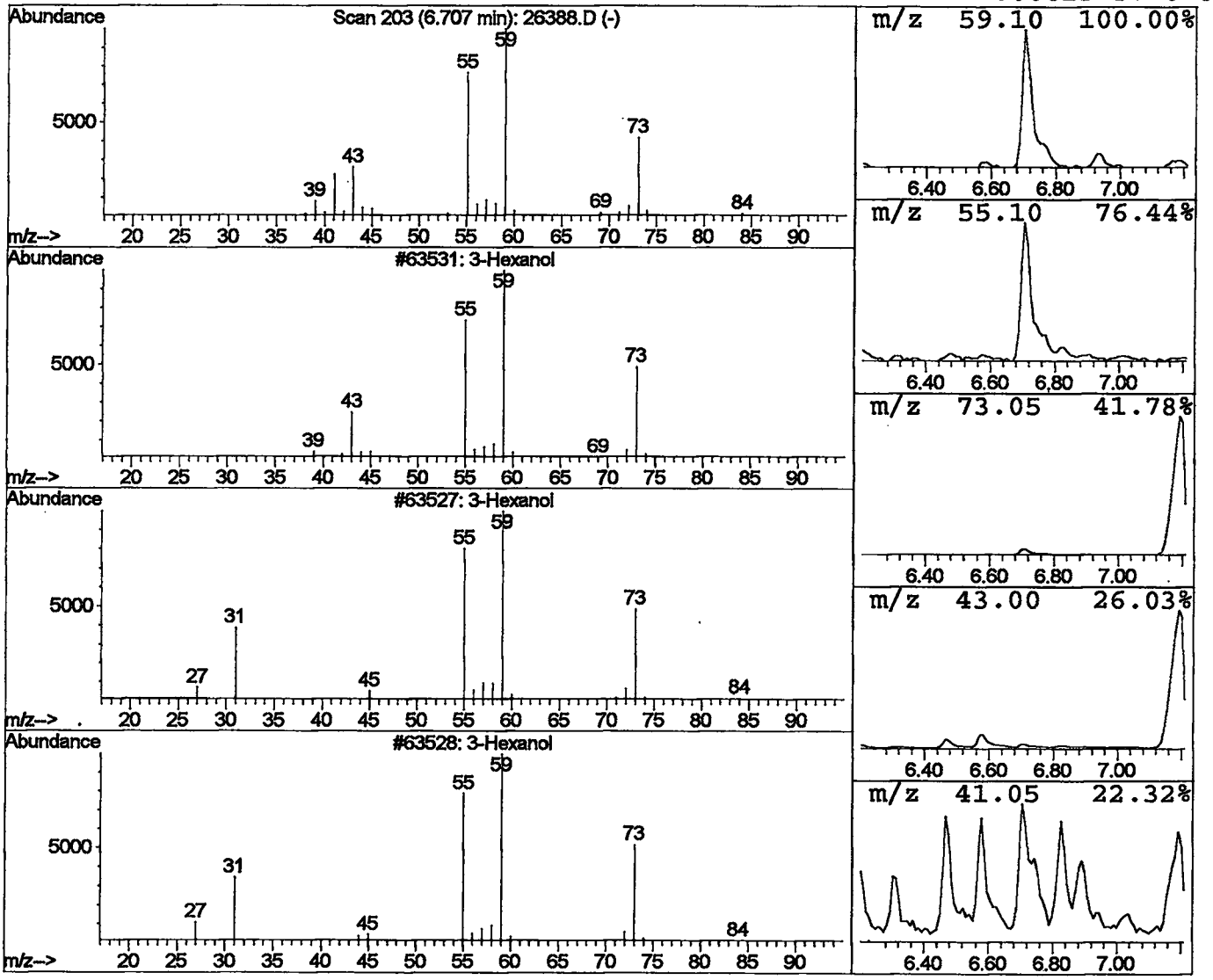
Quant Method : C:\HPCHEM\1\METHODS\NP112701.M (RTE Integrator)
Title : 208 625/TCLP calibration table
Library : C:\DATABASE\NBS75K.L

****

Peak Number 8 3-Hexanol Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.71	0.84 ng/uL	370594	1,4-Dichlorobenzene-d4	11.89

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	3-Hexanol	102	C6H14O	000623-37-0	90
2	3-Hexanol	102	C6H14O	000623-37-0	83
3	3-Hexanol	102	C6H14O	000623-37-0	83
4	3-Hexanol	102	C6H14O	000623-37-0	64



Library Search Compound Report

Data File : C:\HPCHEM\1\DATA\NOV3001\26388.D
Acq On : 1 Dec 2001 12:02 am
Sample : gc/ms unknown 11/3
Misc :
MS Integration Params: LSCINT.P

Vial: 14
Operator: GALOS
Inst : GC/MS Ins
Multiplr: 1.00

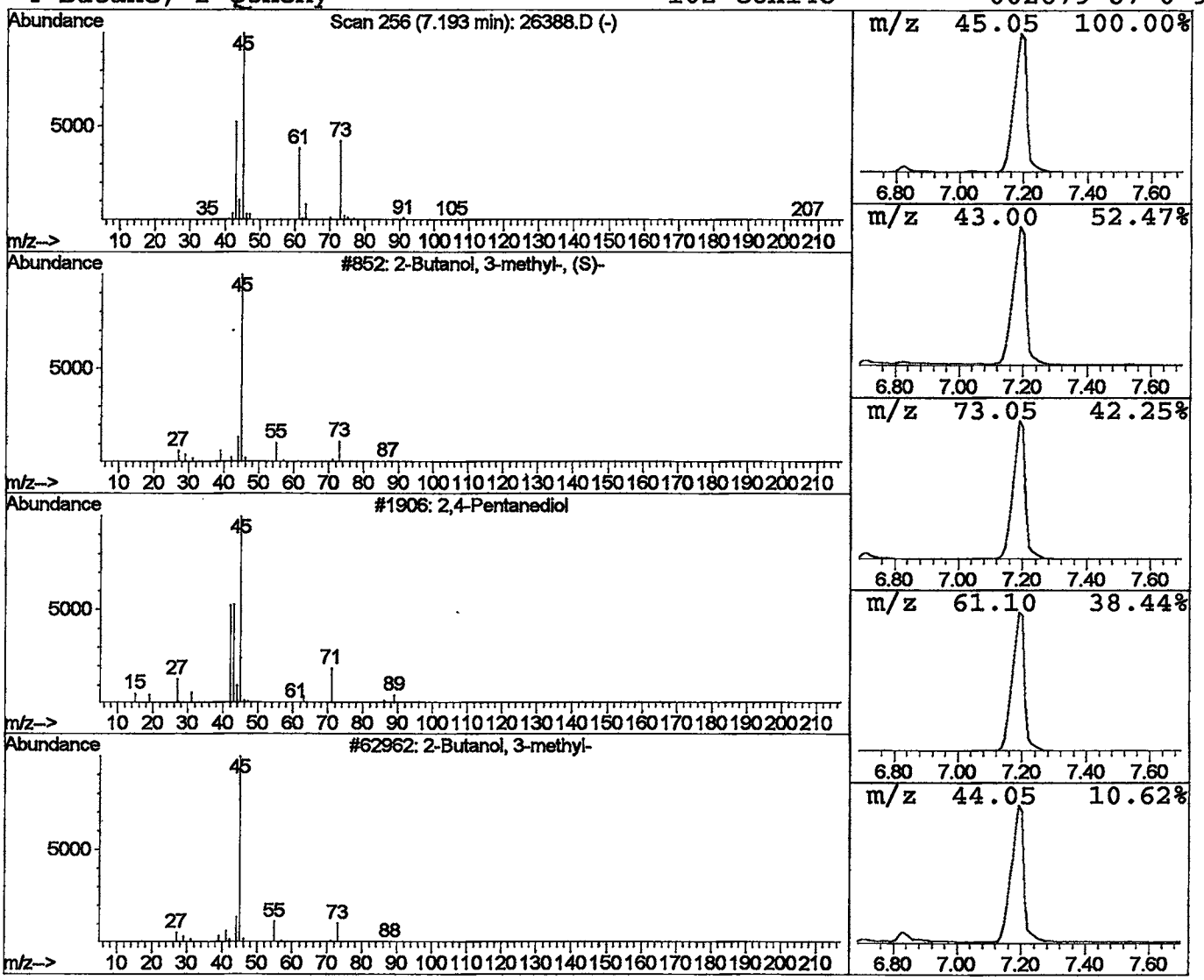
Quant Method : C:\HPCHEM\1\METHODS\NP112701.M (RTE Integrator)
Title : 208 625/TCLP calibration table
Library : C:\DATABASE\NBS75K.L

****

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.19	18.37 ng/uL	8078200	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	Butane, 2-ethoxy-	102	C6H14O	002679-87-0	9



Library Search Compound Report

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

Vial: 14
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Quant Method : C:\HPCHEM\1\METHODS\NP112701.M (RTE Integrator)
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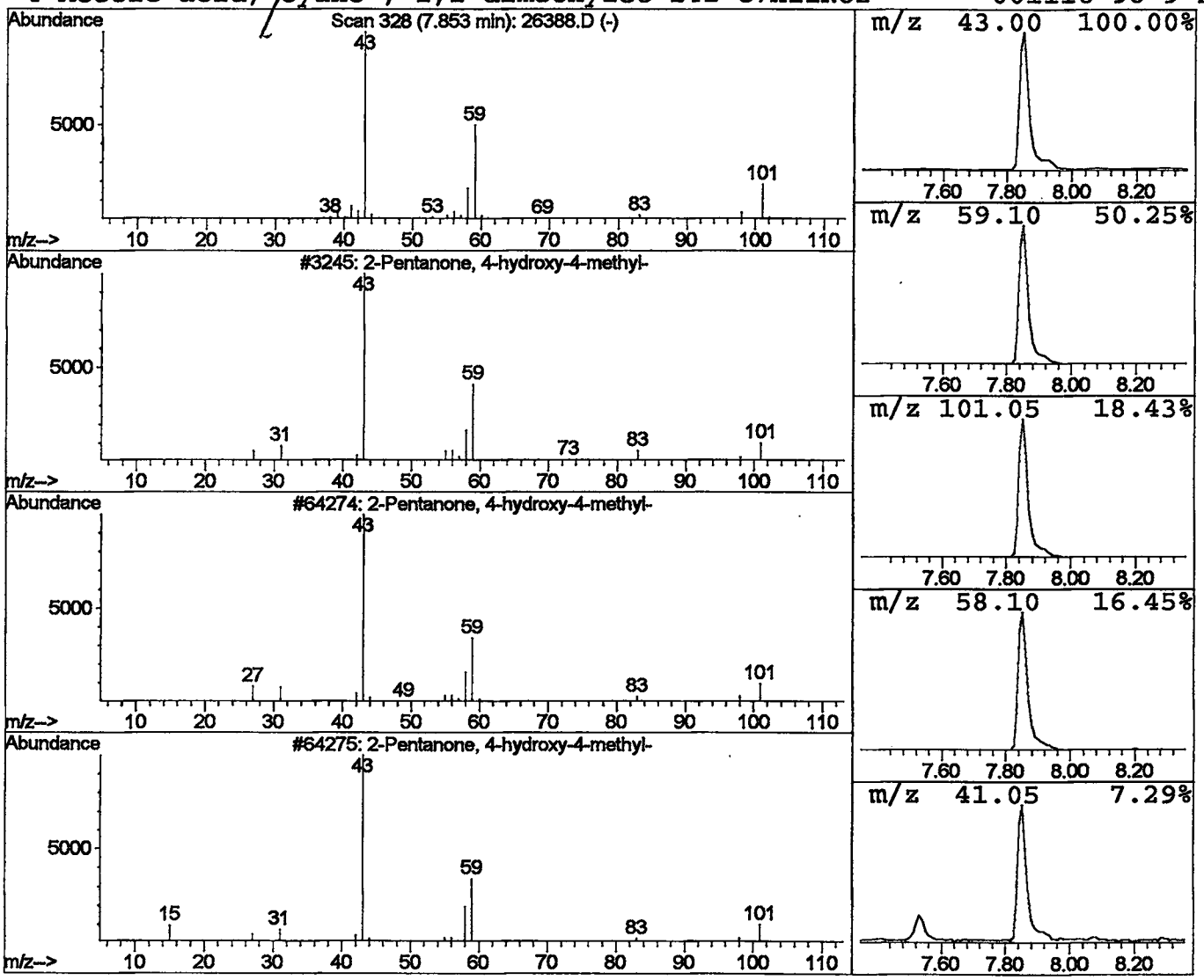
****

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.85	5.35 ng/uL	2350040	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	43
2	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	42
3	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	25
4	Acetic acid, cyano-, 1,1-dimethylet	141	C7H11NO2	001116-98-9	23



Library Search Compound Report

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Acq On : 1 Dec 2001 12:02 am
Sample : gc/ms unknown 11/3
Misc :
MS Integration Params: LSCINT.P

Vial: 14
Operator: GALOS
Inst : GC/MS Ins
Multiplr: 1.00

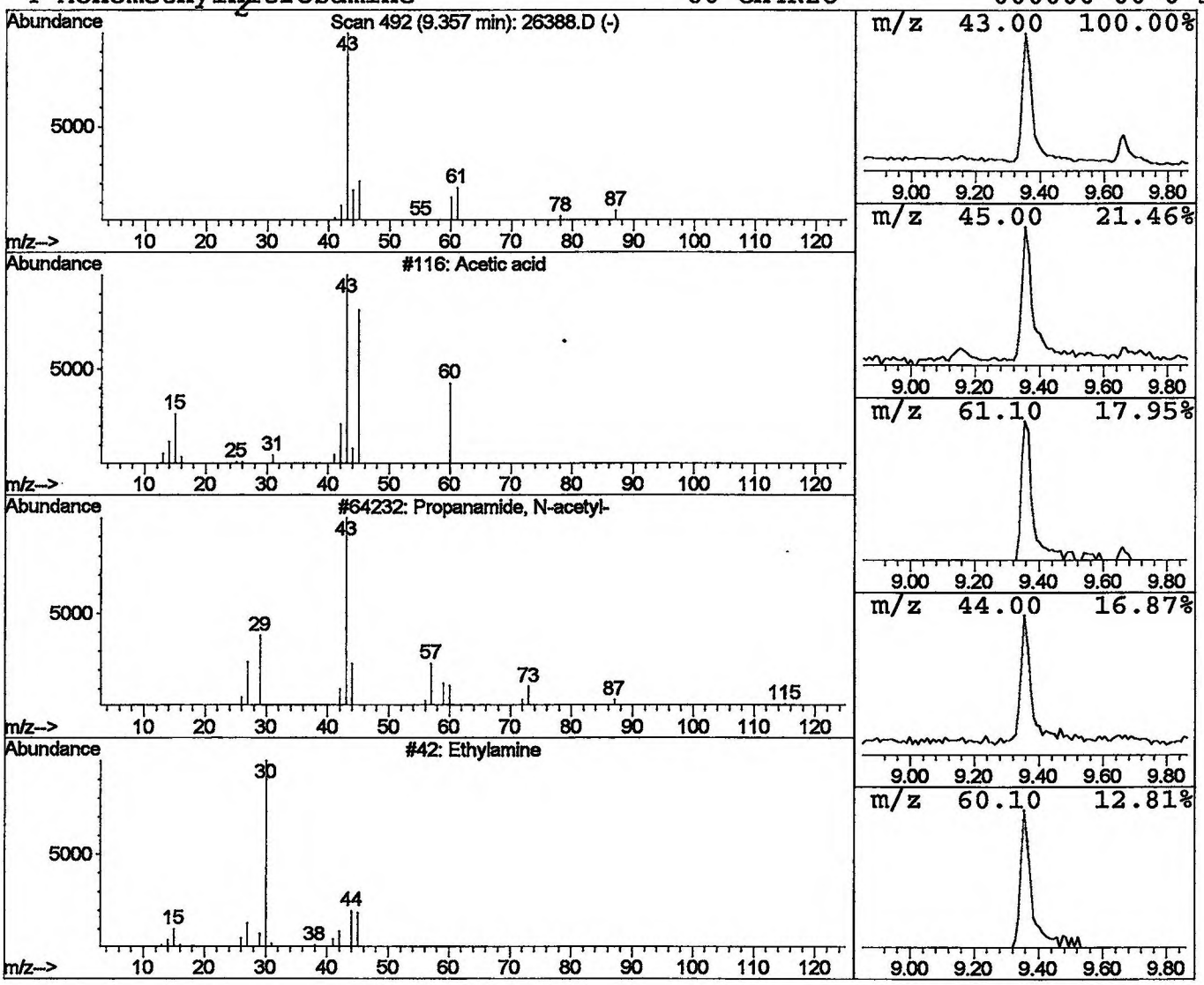
Quant Method : C:\HPCHEM\1\METHODS\NP112701.M (RTE Integrator)
Title : 208 625/TCLP calibration table
Library : C:\DATABASE\NBS75K.L

****

Peak Number 11 Acetic acid — Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.36	0.64 ng/uL	281259	1,4-Dichlorobenzene-d4	11.89

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



Library Search Compound Report

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

Vial: 14
Operator: GALOS
Inst : GC/MS Ins
Multiplr: 1.00

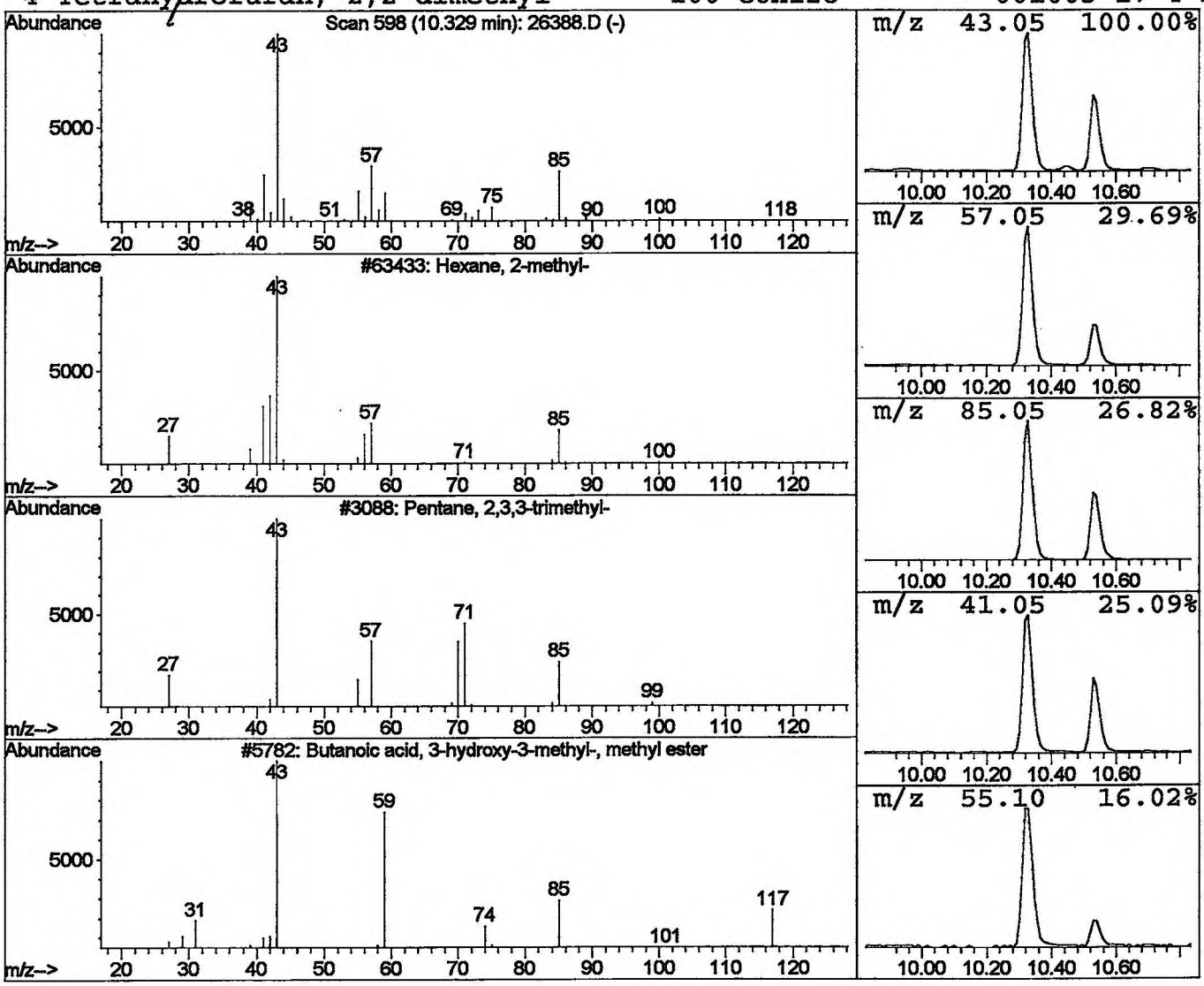
Quant Method : C:\HPCHEM\1\METHODS\NP112701.M (RTE Integrator)
Title : 208 625/TCLP calibration table
Library : C:\DATABASE\NBS75K.L

****

Peak Number 12 Hexane, 2-methyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.33	4.49 ng/uL	1975690	1,4-Dichlorobenzene-d4	11.89

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Hexane, 2-methyl-	100	C7H16	000591-76-4	37
2	Pentane, 2,3,3-trimethyl-	114	C8H18	000560-21-4	37
3	Butanoic acid, 3-hydroxy-3-methyl-,	132	C6H12O3	006149-45-7	28
4	Tetrahydrofuran, 2,2-dimethyl-	100	C6H12O	001003-17-4	25



Library Search Compound Report

Data File : C:\HPCHEM\1\DATA\NOV3001\26388.D
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Misc :
MS Integration Params: LSCINT.P

Vial: 14
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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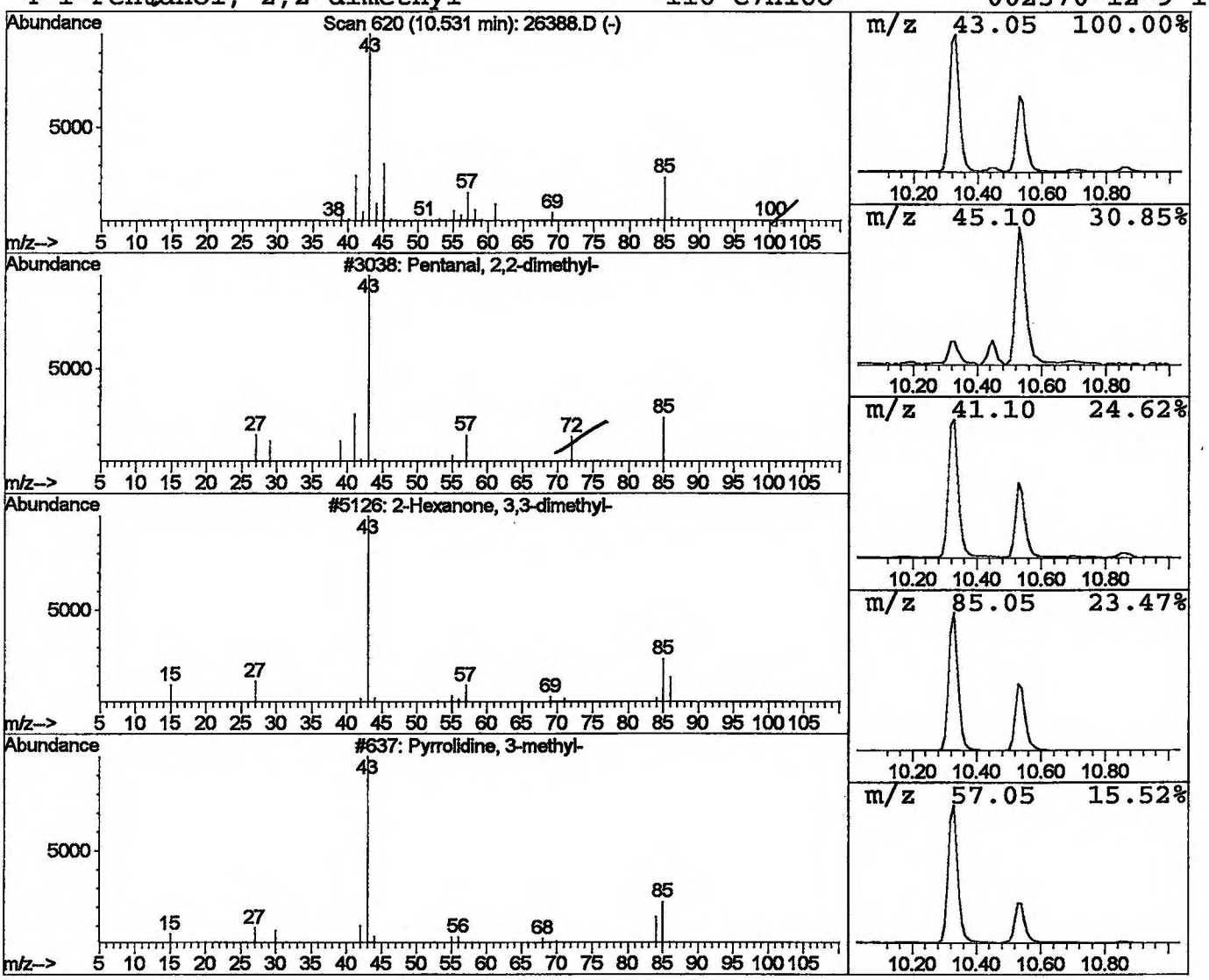
Peak Number 13 Pentanal, 2,2-dimethyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.53	2.15 ng/uL	947159	1,4-Dichlorobenzene-d4	11.89

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Pentanal, 2,2-dimethyl-	114	C7H14O	014250-88-5	37
2	2-Hexanone, 3,3-dimethyl-	128	C8H16O	026118-38-7	32
3	Pyrrolidine, 3-methyl-	85	C5H11N	034375-89-8	32
4	1-Pentanol, 2,2-dimethyl-	116	C7H16O	002370-12-9	17



Library Search Compound Report

Data File : C:\HPCHEM\1\DATA\NOV3001\26388.D
Acq On : 1 Dec 2001 12:02 am
Sample : gc/ms unknown 11/3
Misc :
MS Integration Params: LSCINT.P

Vial: 14
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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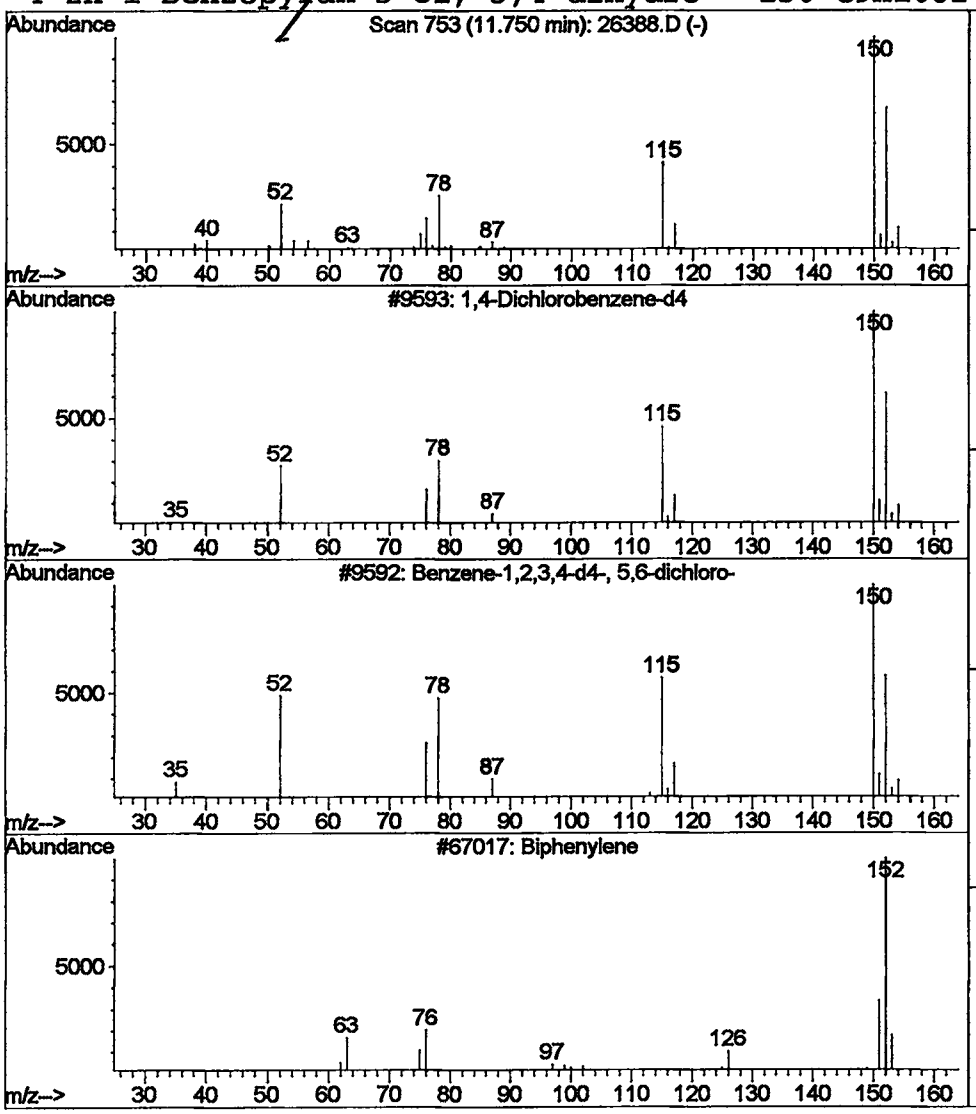
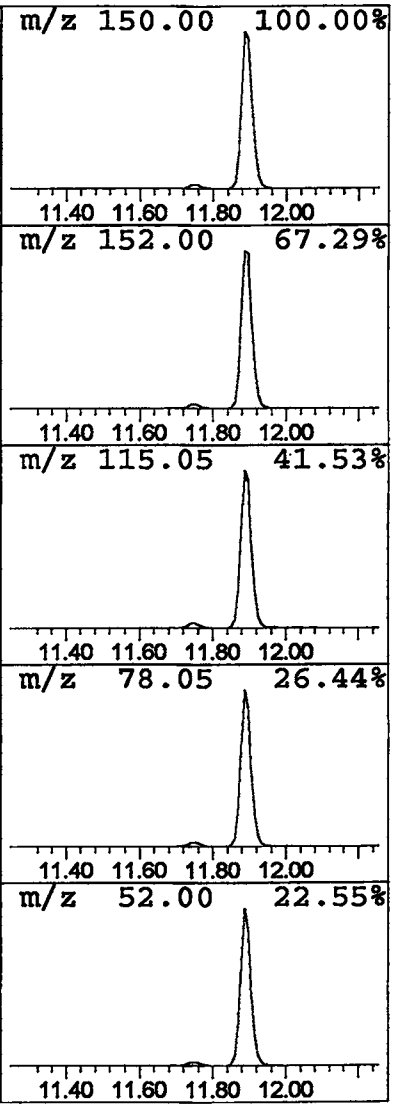
Peak Number 14 1,4-Dichlorobenzene-d4 Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.75	0.57 ng/uL	251345	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	95
2	Benzene-1,2,3,4-d4-, 5,6-dichloro-	150	C6Cl2D4	002199-69-1	42
3	Biphenylene	152	C12H8	000259-79-0	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: 1 Dec 2001 12:02 am
Data File: C:\HPCHEM\1\DATA\NOV3001\26388.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
3-Hexene-2,5-diol	5.02	1.3	ng/uL	591056	ISTD01	11.89	8792770	20.0
1-Butanol	5.12	2.0	ng/uL	865260	ISTD01	11.89	8792770	20.0
Oxirane, 3-ethyl-2,2	5.18	4.0	ng/uL	1753570	ISTD01	11.89	8792770	20.0
Oxirane, 2-methyl-3-	5.67	0.6	ng/uL	270347	ISTD01	11.89	8792770	20.0
2-Furanmethanamine,	6.12	2.3	ng/uL	1027330	ISTD01	11.89	8792770	20.0
3-Hexanone	6.47	0.6	ng/uL	256746	ISTD01	11.89	8792770	20.0
2-Hexanone	6.58	0.7	ng/uL	304613	ISTD01	11.89	8792770	20.0
3-Hexanol	6.71	0.8	ng/uL	370594	ISTD01	11.89	8792770	20.0
2-Butanol, 3-methyl-	7.19	18.4	ng/uL	8078200	ISTD01	11.89	8792770	20.0
2-Pentanone, 4-hydro	7.85	5.3	ng/uL	2350040	ISTD01	11.89	8792770	20.0
Acetic acid	9.36	0.6	ng/uL	281259	ISTD01	11.89	8792770	20.0
Hexane, 2-methyl-	10.33	4.5	ng/uL	1975690	ISTD01	11.89	8792770	20.0
Pentanal, 2,2-dimeth	10.53	2.2	ng/uL	947159	ISTD01	11.89	8792770	20.0
1,4-Dichlorobenzene-	11.75	0.6	ng/uL	251345	ISTD01	11.89	8792770	20.0

26388.D NP112701.M Tue Dec 04 10:40:28 2001