Quantitation Report (QT Reviewed)

Data File : C:\HPCHEM\1\DATA\DEC2101\LB.D
Acq On : 22 Dec 2001 2:07 am
Sample : gcms unknown 11/23
Misc :
28169-28176
28104-28106
Vial: 10
Operator: GALOS
Inst : GC/MS Ins
Multiplr: 1.00
MS Integration Params: RTEINT.P
Quant Time: Dec 28 11:50 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	11.80	152	691607	20.00	ng/uL	-0.14
19) Naphthalene-d8	15.41	136	2704570	20.00	ng/uL	-0.14
31) Acenaphthene-d10	20.67	164	1191938	20.00	ng/uL	-0.15
49) Phenanthrene-d10	25.08	188	1736568m	20.00	ng/uL	-0.16
59) Chrysene-d12	33.09	240	1178288	20.00	ng/uL	-0.16
68) Perylene-d12	37.15	264	713135	20.00	ng/uL	-0.19
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	11.80	132	317	0.01	ng/uL	0.37
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	18.83	172	1187	0.02	ng/uL	0.00
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	0.00	45	0	N.D.		
16) 3&4-Methylphenol	0.00	108	0	N.D.		

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

Quantitation Report (QT Reviewed)

Data File : C:\HPCHEM\1\DATA\DEC2101\LB.D Vial: 10
Acq On : 22 Dec 2001 2:07 am Operator: GALOS
Sample : gcms unknown 11/23 Inst : GC/MS Ins
Misc : Multiplr: 1.00
MS Integration Params: RTEINT.P
Quant Time: Dec 28 11:50 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

Compound	R.T.	QIon	Response	Conc Unit	Qvalue
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.	
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.	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	27.07	149	1394	N.D.	
58) Fluoranthene	0.00	202	0	N.D.	
60) 3,3'-Dichlorobenzidine	0.00	252	0	N.D.	

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

Quantitation Report (QT Reviewed)

Data File : C:\HPCHEM\1\DATA\DEC2101\LB.D Vial: 10
Acq On : 22 Dec 2001 2:07 am Operator: GALOS
Sample : gcms unknown 11/23 Inst : GC/MS Ins
Misc : Multiplr: 1.00
MS Integration Params: RTEINT.P
Quant Time: Dec 28 11:50 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

	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	33.09	228	2822	N.D.	
66)	Bis(2-ethylhexyl)phthalate	33.36	149	8372	N.D.	
67)	Chrysene	33.09	228	2822	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	37.15	252	2450	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.	

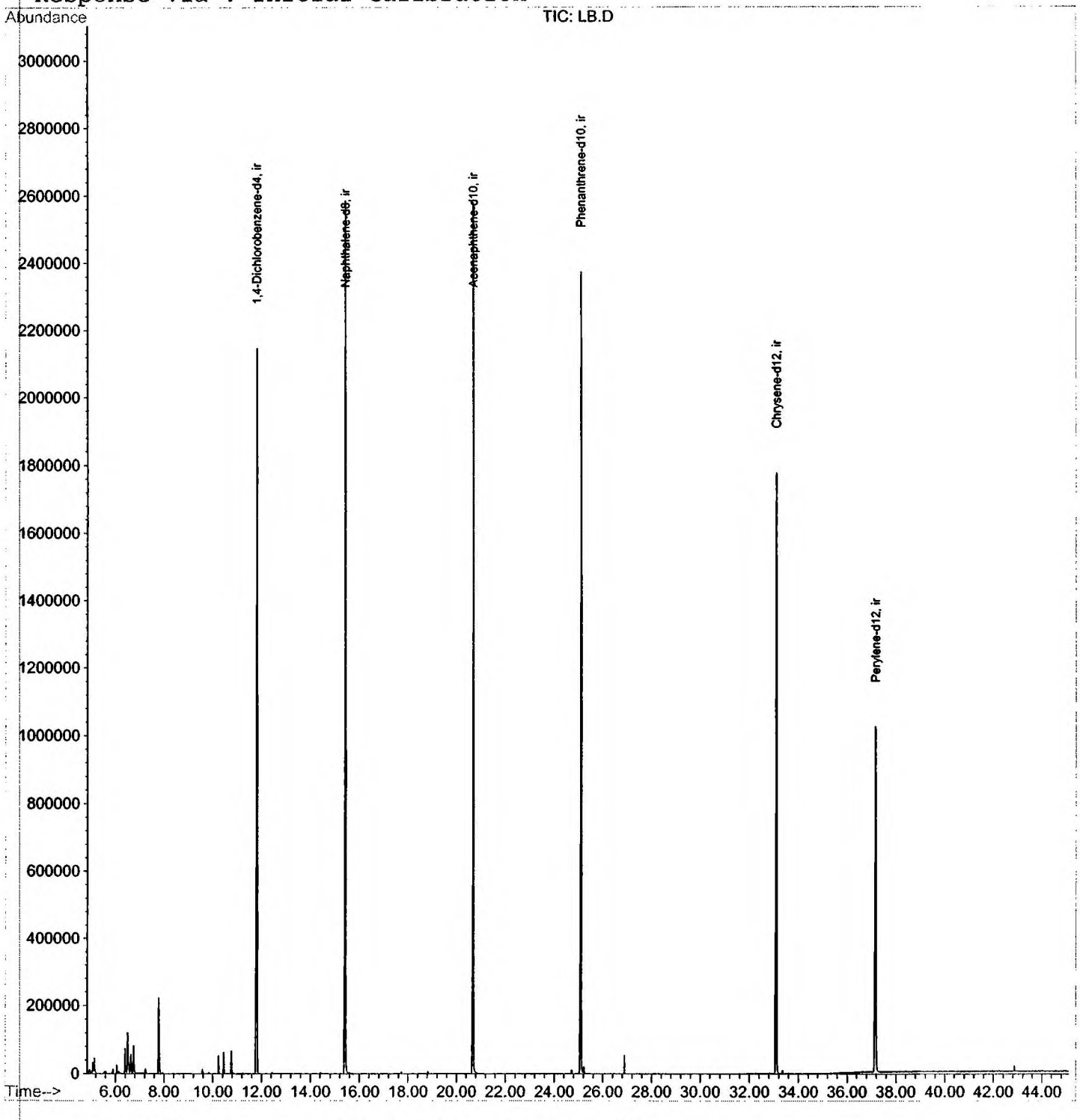
(#) = qualifier out of range (m) = manual integration

Quantitation Report

Data File : C:\HPCHEM\1\DATA\DEC2101\LB.D
Acq On : 22 Dec 2001 2:07 am
Sample : gcms unknown 11/23
Misc :
MS Integration Params: RTEINT.P
Quant Time: Dec 28 11:50 2001

Vial: 10
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\DEC2101\LB.D Vial: 10
Acq On : 22 Dec 2001 2:07 am Operator: GALOS
Sample : gcms unknown 11/23 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

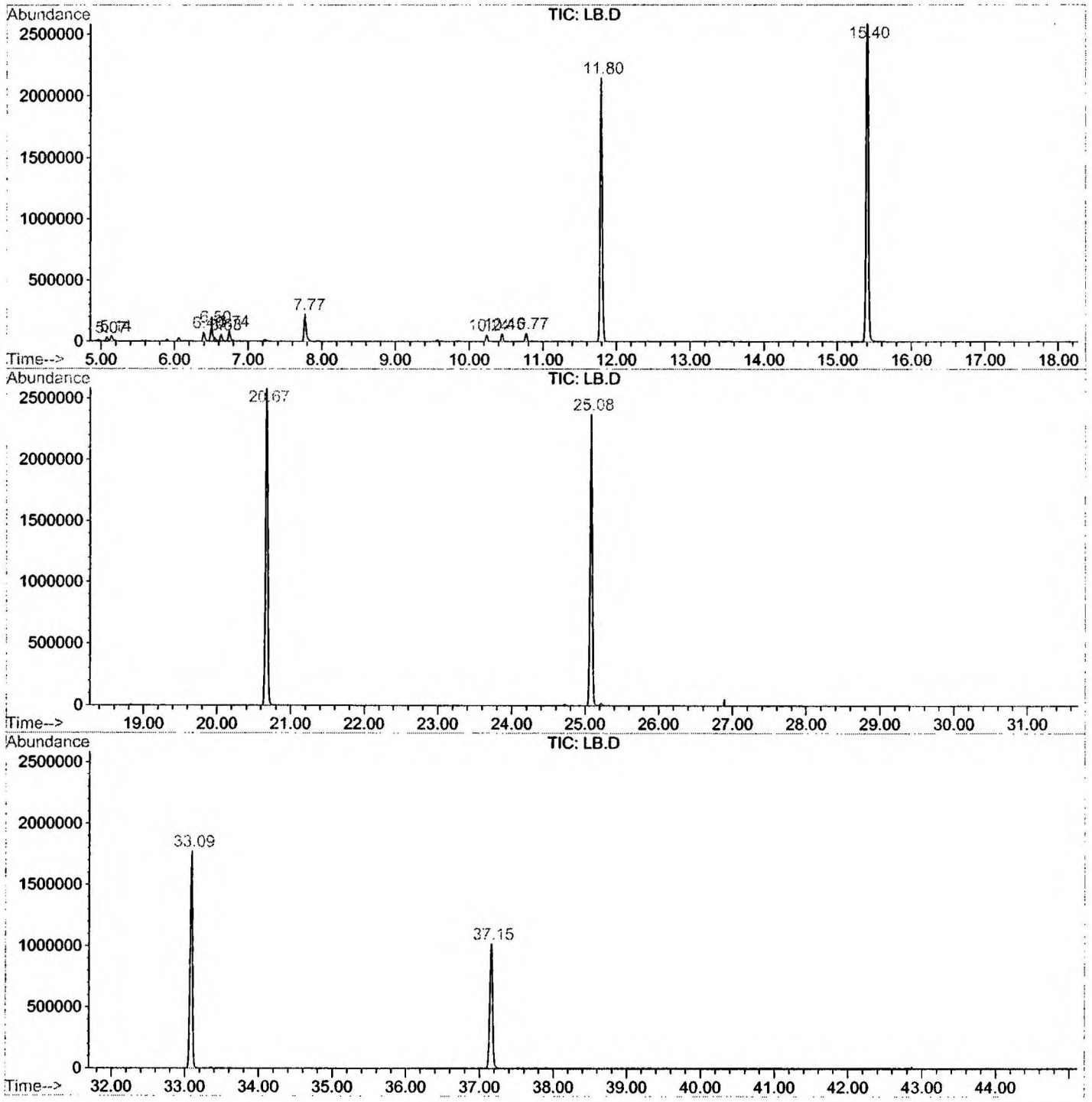
peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	peak area	peak % max.	% of total
1	5.075	22	25	29	rVV	33539	61671	1.11%	0.217%
2	5.139	29	32	41	rVB	46409	98014	1.77%	0.345%
3	6.395	165	169	176	rBV	74313	144116	2.60%	0.507%
4	6.496	176	180	184	rBV	118215	242562	4.38%	0.853%
5	6.634	191	195	203	rVB	55653	107868	1.95%	0.380%
6	6.744	203	207	217	rBV	81706	163024	2.94%	0.574%
7	7.771	315	319	332	rBB	223564	448535	8.10%	1.578%
8	10.237	584	588	593	rBV	53525	95992	1.73%	0.338%
9	10.448	607	611	616	rBB	63833	116114	2.10%	0.409%
10	10.769	642	646	651	rVB	67743	120862	2.18%	0.425%
11	11.796	752	758	769	rVB	2148170	4319372	78.03%	15.198%
12	15.400	1145	1151	1161	rBV	2586279	5535834	100.00%	19.478%
13	20.673	1720	1726	1742	rBV	2580649	5346719	96.58%	18.812%
14	25.084	2200	2207	2218	rBV2	2374517	4963957	89.67%	17.466%
15	33.089	3073	3080	3091	rBV2	1777154	3877708	70.05%	13.644%
16	37.152	3515	3523	3536	rBV2	1020636	2778780	50.20%	9.777%

Sum of corrected areas: 28421128

LB.D NP112701.M Fri Dec 28 11:46:36 2001

LSC Report - Integrated Chromatogram

File : C:\HPCHEM\1\DATA\DEC2101\LB.D
Operator : GALOS
Acquired : 22 Dec 2001 2:07 am using AcqMethod NP112701
Instrument : GC/MS Ins
Sample Name: gcms unknown 11/23
Misc Info :
Vial Number: 10
Quant File :NP112701.RES (RTE Integrator)



LB.D NP112701.M Fri Dec 28 11:46:39 2001

Library Search Compound Report

Data File : C:\HPCHEM\1\DATA\DEC2101\LB.D
Acq On : 22 Dec 2001 2:07 am
Sample : gcms unknown 11/23
Misc :
MS Integration Params: LSCINT.P

Vial: 10
Operator: GALOS
Inst : GC/MS Ins
Multiplr: 1.00

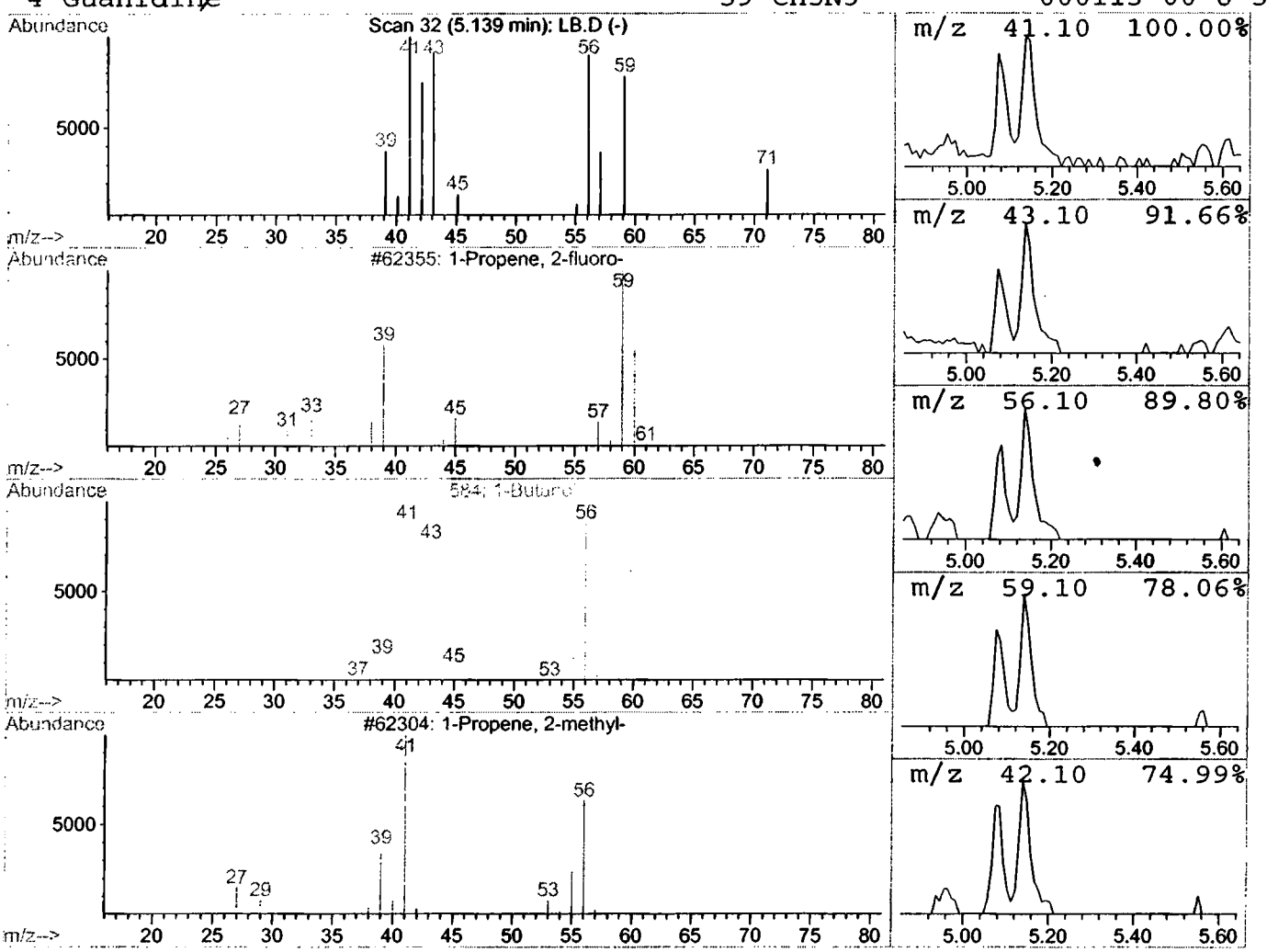
Quant Method : C:\HPCHEM\1\METHODS\NP112701.M (RTE Integrator)
Title : 208 625/TCLP calibration table
Library : C:\DATABASE\NBS75K.L

****

Peak Number 1 1-Propene, 2-fluoro- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.14	0.45 ng/uL	98014	1,4-Dichlorobenzene-d4	11.80

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Propene, 2-fluoro-	60	C3H5F	001184-60-7	9
2	1-Butanol	74	C4H10O	000071-36-3	9
3	1-Propene, 2-methyl-	56	C4H8	000115-11-7	7
4	Guanidine	59	CH5N3	000113-00-8	5



Library Search Compound Report

Data File : C:\HPCHEM\1\DATA\DEC2101\LB.D
Acq On : 22 Dec 2001 2:07 am
Sample : gcms unknown 11/23
Misc :
MS Integration Params: LSCINT.P

Vial: 10
Operator: GALOS
Inst : GC/MS Ins
Multiplr: 1.00

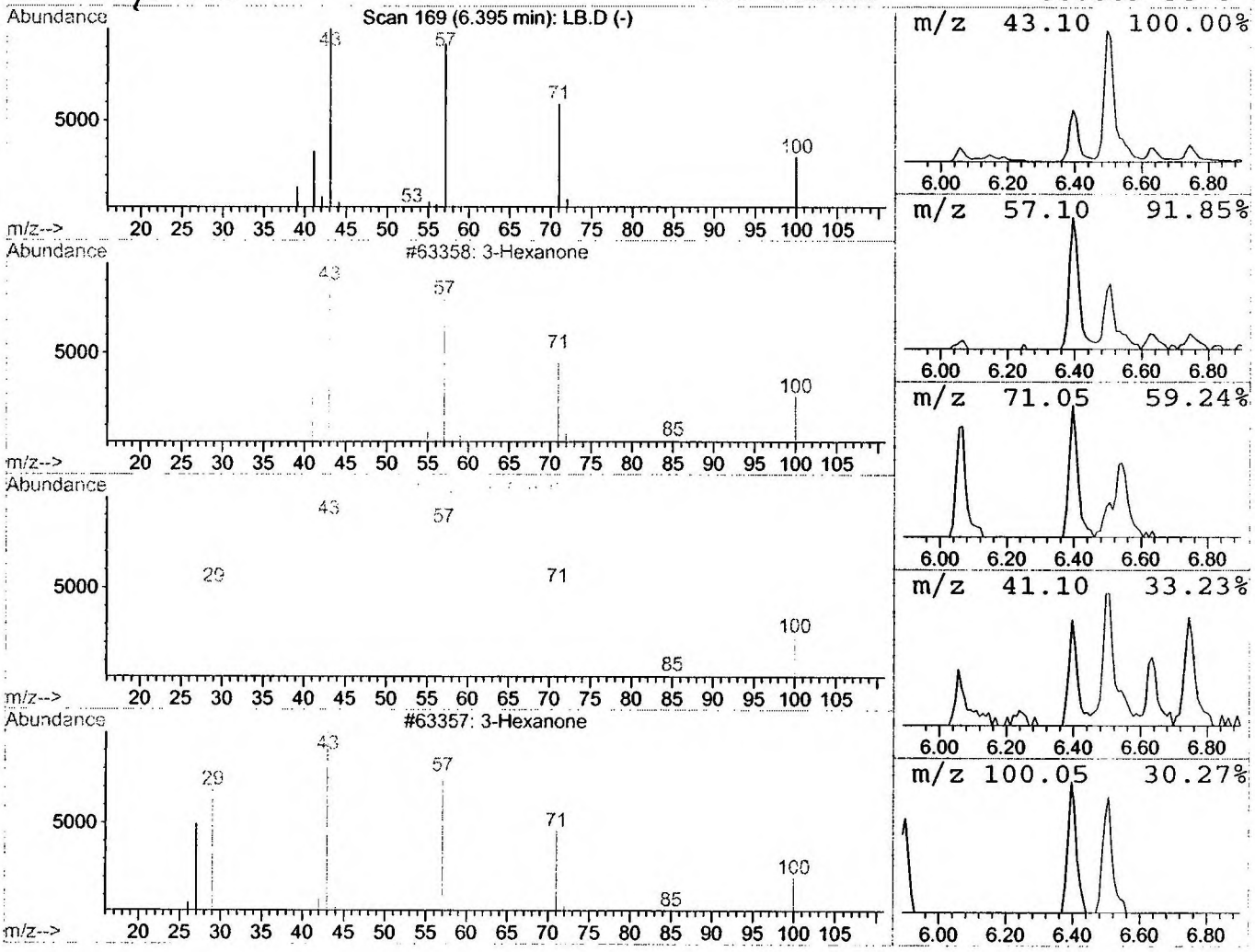
Quant Method : C:\HPCHEM\1\METHODS\NP112701.M (RTE Integrator)
Title : 208 625/TCLP calibration table
Library : C:\DATABASE\NBS75K.L

****

Peak Number 2 3-Hexanone Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.40	0.67 ng/uL	144116	1,4-Dichlorobenzene-d4	11.80

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	3-Hexanone	100	C6H12O	000589-38-8	91
2	3-Hexanone	100	C6H12O	000589-38-8	90
3	3-Hexanone	100	C6H12O	000589-38-8	83
4	3-Hexanone	100	C6H12O	000589-38-8	56



Library Search Compound Report

Data File : C:\HPCHEM\1\DATA\DEC2101\LB.D
Acq On : 22 Dec 2001 2:07 am
Sample : gcms unknown 11/23
Misc :
MS Integration Params: LSCINT.P

Vial: 10
Operator: GALOS
Inst : GC/MS Ins
Multiplr: 1.00

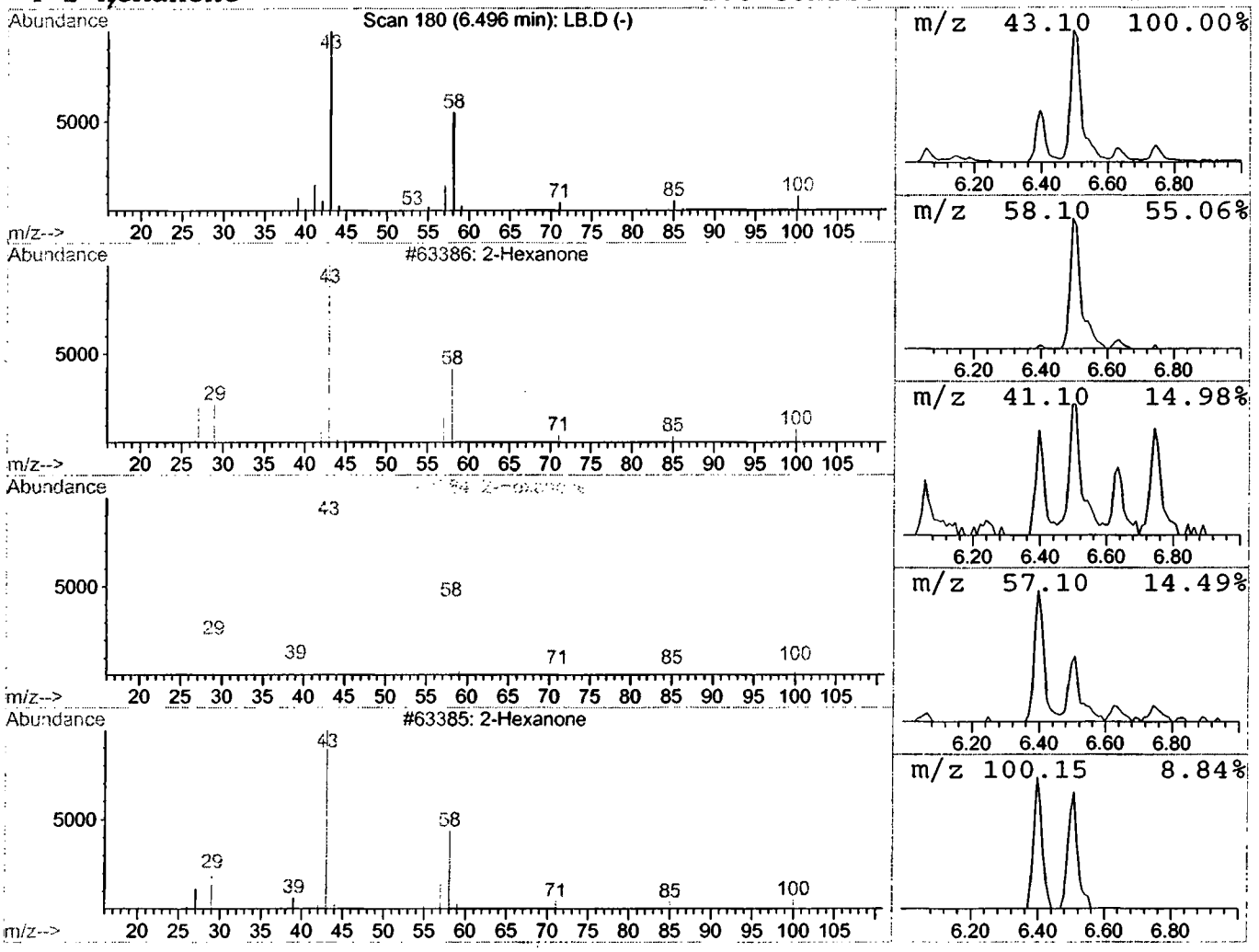
Quant Method : C:\HPCHEM\1\METHODS\NP112701.M (RTE Integrator)
Title : 208 625/TCLP calibration table
Library : C:\DATABASE\NBS75K.L

****

Peak Number 3 2-Hexanone Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.50	1.12 ng/uL	242562	1,4-Dichlorobenzene-d4	11.80

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Hexanone	100	C6H12O	000591-78-6	91
2	2-Hexanone	100	C6H12O	000591-78-6	91
3	2-Hexanone	100	C6H12O	000591-78-6	91
4	2-Hexanone	100	C6H12O	000591-78-6	90



Library Search Compound Report

Data File : C:\HPCHEM\1\DATA\DEC2101\LB.D Vial: 10
Acq On : 22 Dec 2001 2:07 am Operator: GALOS
Sample : gcms unknown 11/23 Inst : GC/MS Ins
Misc : Multiplr: 1.00
MS Integration Params: LSCINT.P

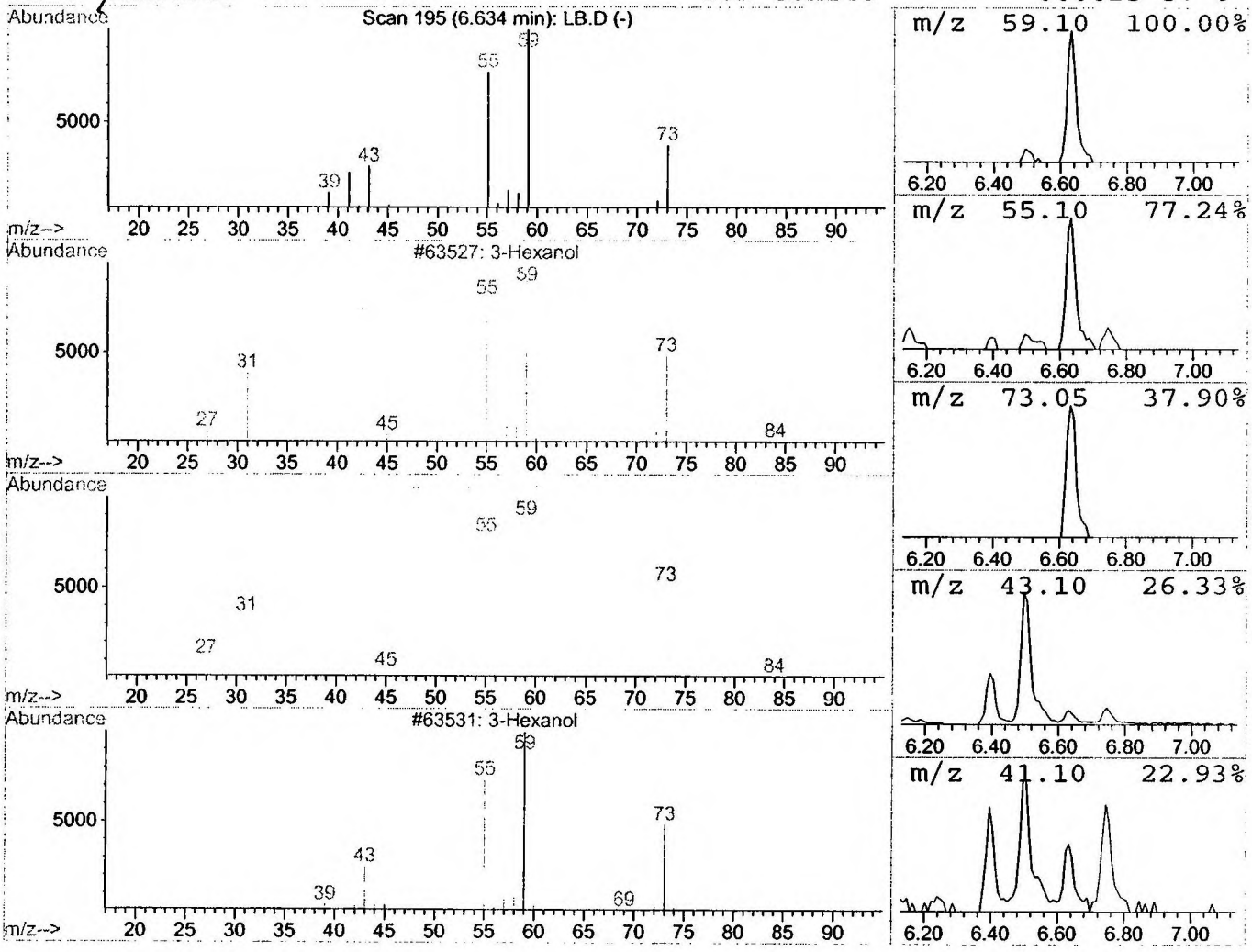
Quant Method : C:\HPCHEM\1\METHODS\NP112701.M (RTE Integrator)
Title : 208 625/TCLP calibration table
Library : C:\DATABASE\NBS75K.L

****

Peak Number 4 3-Hexanol Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.63	0.50 ng/uL	107868	1,4-Dichlorobenzene-d4	11.80

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	83
4	3-Hexanol	102	C6H14O	000623-37-0	64



Library Search Compound Report

Data File : C:\HPCHEM\1\DATA\DEC2101\LB.D
Acq On : 22 Dec 2001 2:07 am
Sample : gcms unknown 11/23
Misc :
MS Integration Params: LSCINT.P

Vial: 10
Operator: GALOS
Inst : GC/MS Ins
Multiplr: 1.00

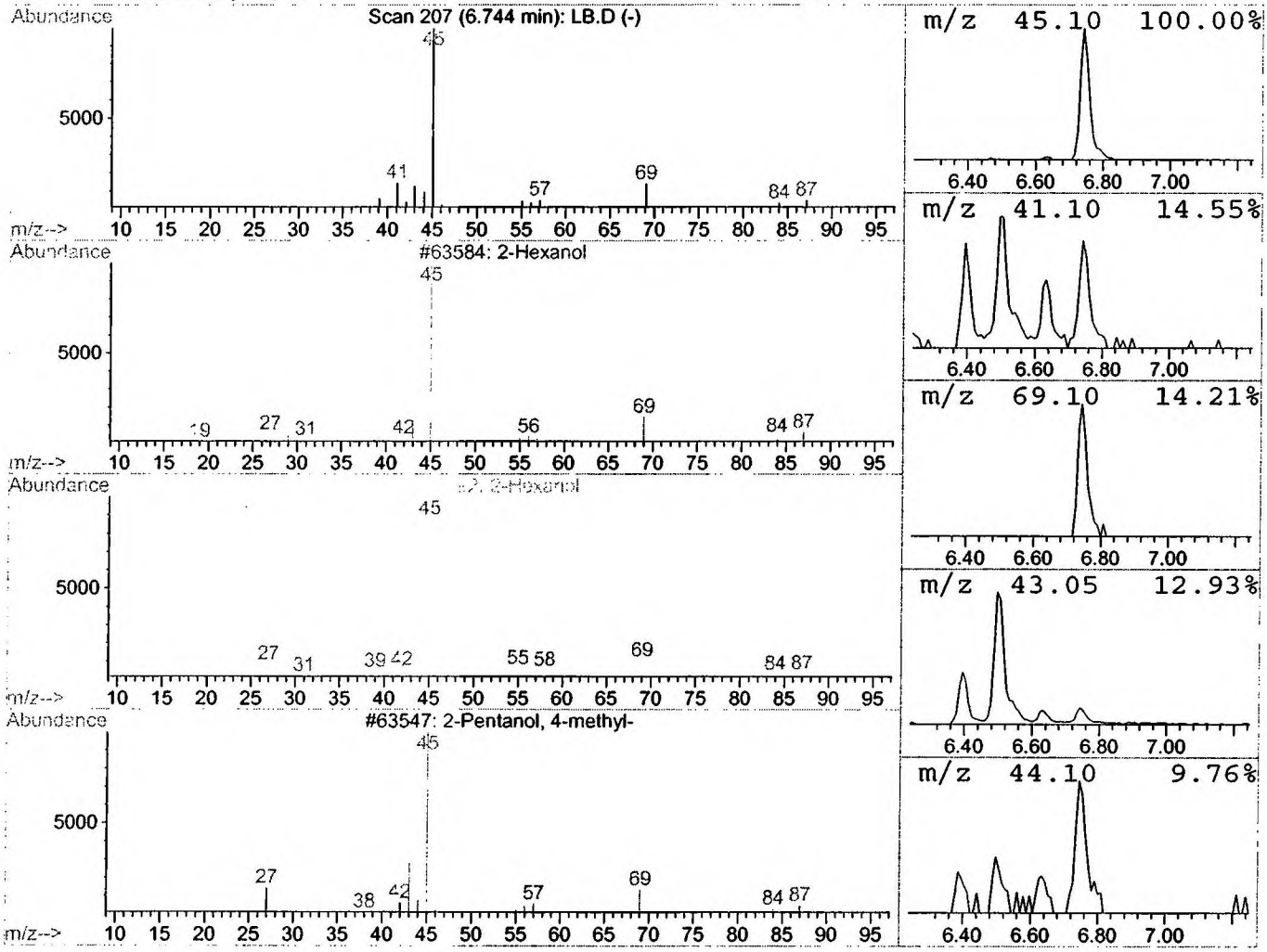
Quant Method : C:\HPCHEM\1\METHODS\NP112701.M (RTE Integrator)
Title : 208 625/TCLP calibration table
Library : C:\DATABASE\NBS75K.L

****

Peak Number 5 2-Hexanol Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.74	0.75 ng/uL	163024	1,4-Dichlorobenzene-d4	11.80

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Hexanol	102	C6H14O	000626-93-7	83
2	2-Hexanol	102	C6H14O	000626-93-7	83
3	2-Pentanol, 4-methyl-	102	C6H14O	000108-11-2	74
4	2-Hexanol, (R)-	102	C6H14O	026549-24-6	74



Library Search Compound Report

Data File : C:\HPCHEM\1\DATA\DEC2101\LB.D
Acq On : 22 Dec 2001 2:07 am
Sample : gcms unknown 11/23
Misc :
MS Integration Params: LSCINT.P

Vial: 10
Operator: GALOS
Inst : GC/MS Ins
Multiplr: 1.00

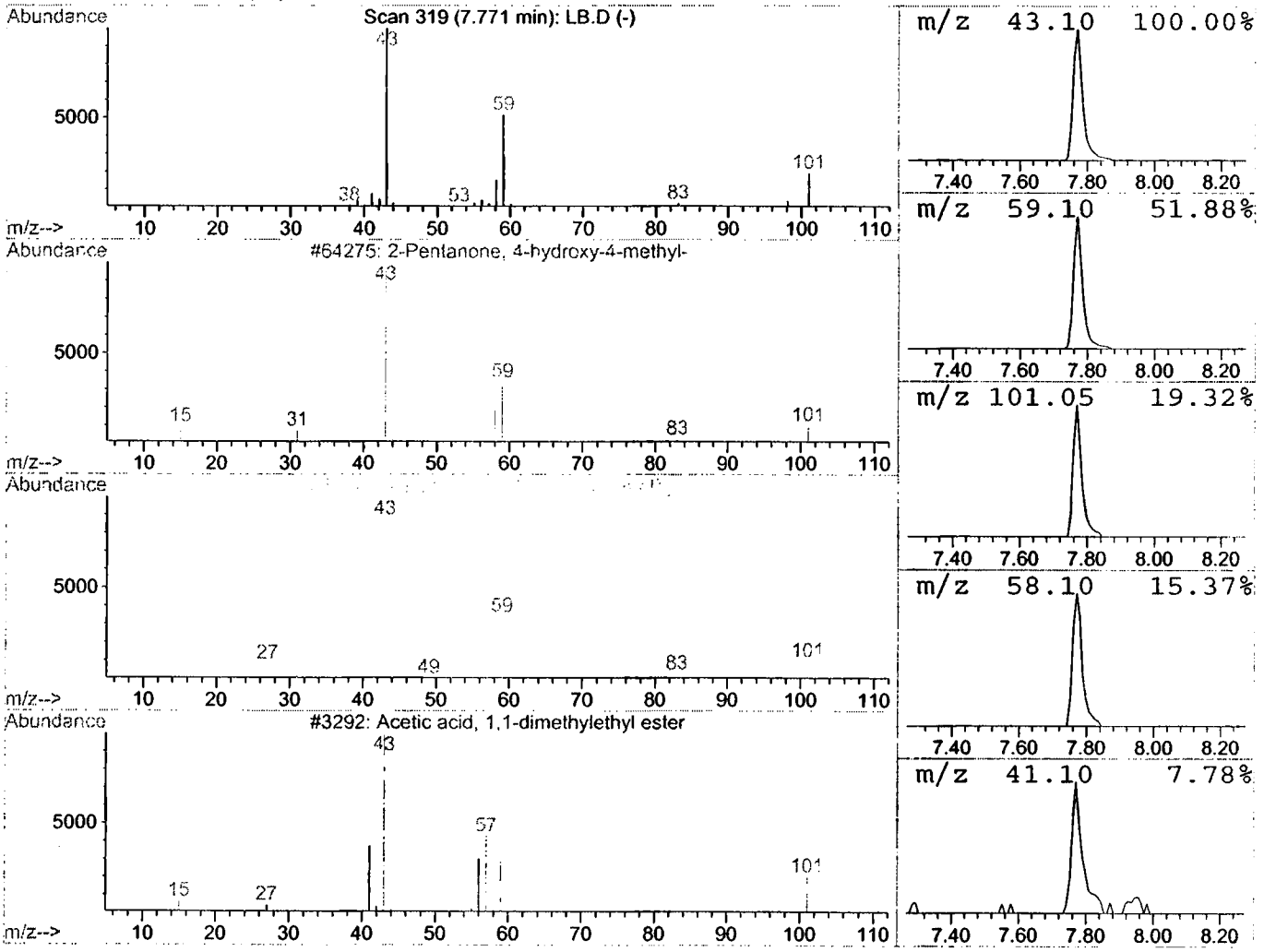
Quant Method : C:\HPCHEM\1\METHODS\NP112701.M (RTE Integrator)
Title : 208 625/TCLP calibration table
Library : C:\DATABASE\NBS75K.L

****

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.77	2.08 ng/uL	448535	1,4-Dichlorobenzene-d4	11.80

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	40
2	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	40
3	Acetic acid, 1,1-dimethylethyl este	116	C6H12O2	000540-88-5	33
4	CH2=CHCH2C(O)OCH3	100	C5H8O2	003724-55-8	10



Library Search Compound Report

Data File : C:\HPCHEM\1\DATA\DEC2101\LB.D
Acq On : 22 Dec 2001 2:07 am
Sample : gcms unknown 11/23
Misc :
MS Integration Params: LSCINT.P

Vial: 10
Operator: GALOS
Inst : GC/MS Ins
Multiplr: 1.00

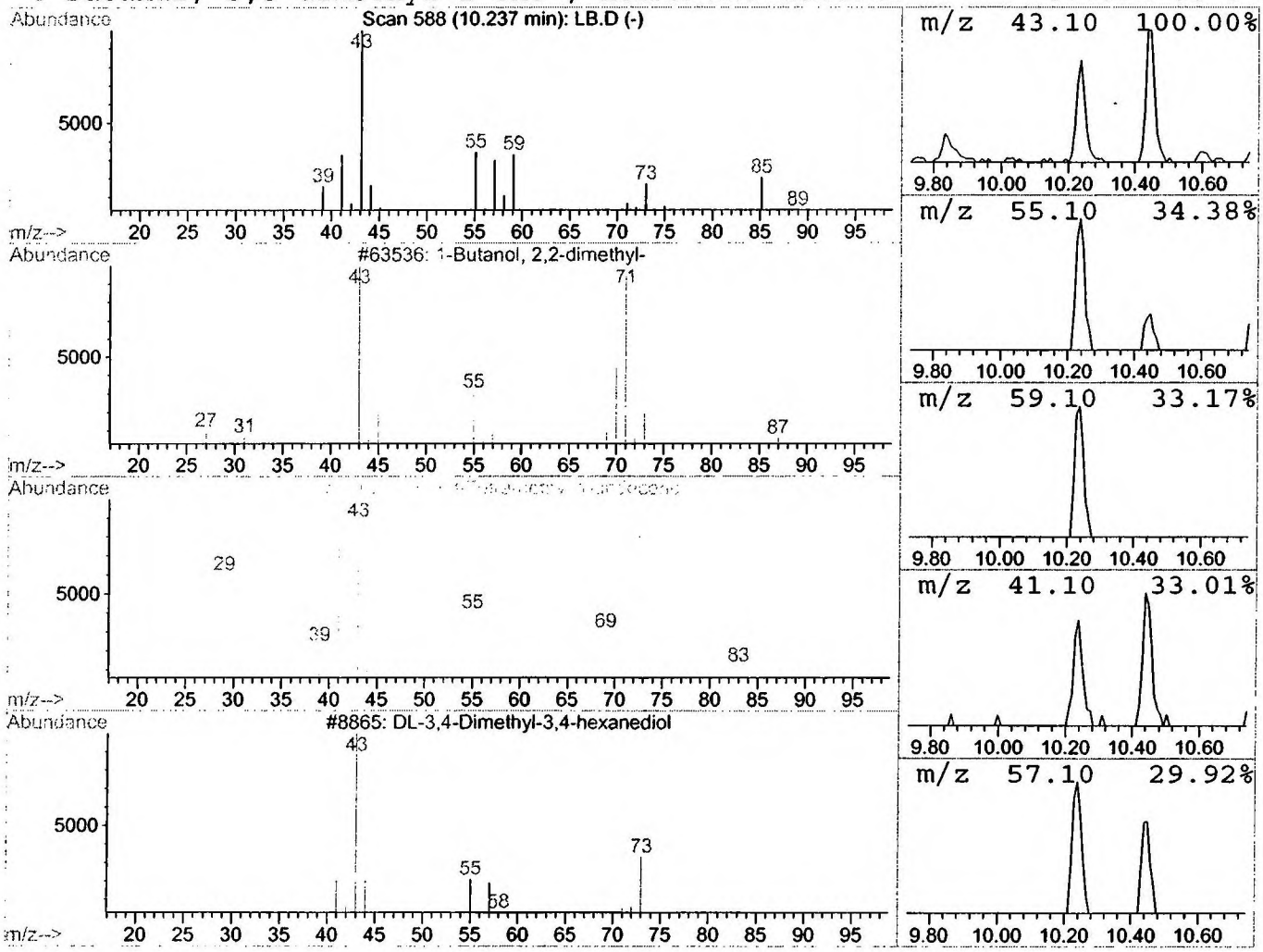
Quant Method : C:\HPCHEM\1\METHODS\NP112701.M (RTE Integrator)
Title : 208 625/TCLP calibration table
Library : C:\DATABASE\NBS75K.L

****

Peak Number 7 1-Butanol, 2,2-dimethyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.24	0.44 ng/uL	95992	1,4-Dichlorobenzene-d4	11.80

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Butanol, 2,2-dimethyl-	102	C6H14O	001185-33-7	9
2	2,4,6,8-Tetramethyl-1-undecene	210	C15H30	059920-26-2	9
3	DL-3,4-Dimethyl-3,4-hexanediol	146	C8H18O2	032388-94-6	9
4	Butanal, 3,3-dimethyl-2-oxo-, hemih	114	C6H10O2	077572-68-0	9



Library Search Compound Report

Data File : C:\HPCHEM\1\DATA\DEC2101\LB.D
Acq On : 22 Dec 2001 2:07 am
Sample : gcms unknown 11/23
Misc :
MS Integration Params: LSCINT.P

Vial: 10
Operator: GALOS
Inst : GC/MS Ins
Multiplr: 1.00

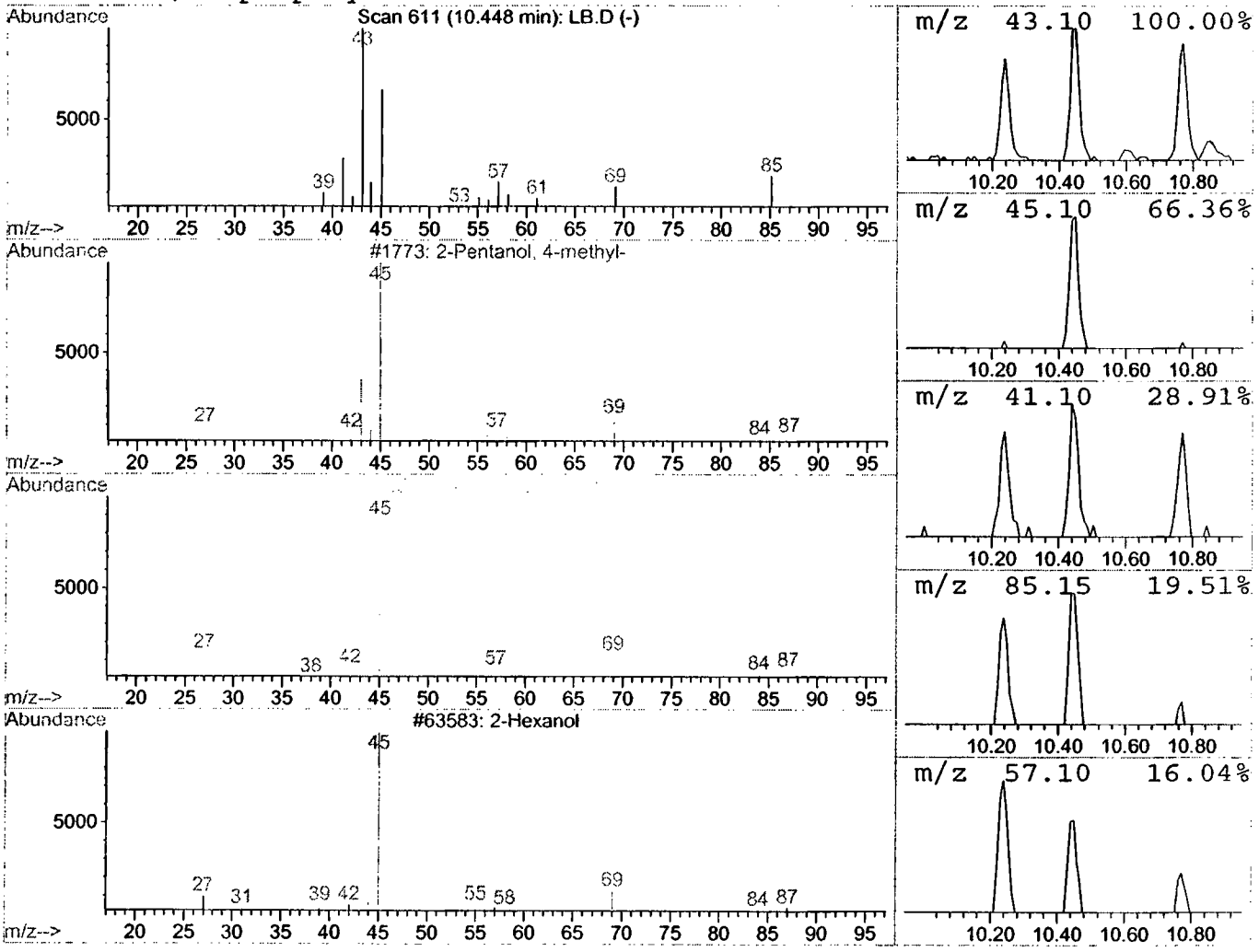
Quant Method : C:\HPCHEM\1\METHODS\NP112701.M (RTE Integrator)
Title : 208 625/TCLP calibration table
Library : C:\DATABASE\NBS75K.L

****

Peak Number 8 2-Pentanol, 4-methyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.45	0.54 ng/uL	116114	1,4-Dichlorobenzene-d4	11.80

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Pentanol, 4-methyl-	102	C6H14O	000108-11-2	25
2	2-Pentanol, 4-methyl-	102	C6H14O	000108-11-2	23
3	2-Hexanol	102	C6H14O	000626-93-7	23
4	Hexane, 1-propoxy-	144	C9H20O	053685-78-2	17



Library Search Compound Report

Data File : C:\HPCHEM\1\DATA\DEC2101\LB.D
Acq On : 22 Dec 2001 2:07 am
Sample : gcms unknown 11/23
Misc :
MS Integration Params: LSCINT.P

Vial: 10
Operator: GALOS
Inst : GC/MS Ins
Multiplr: 1.00

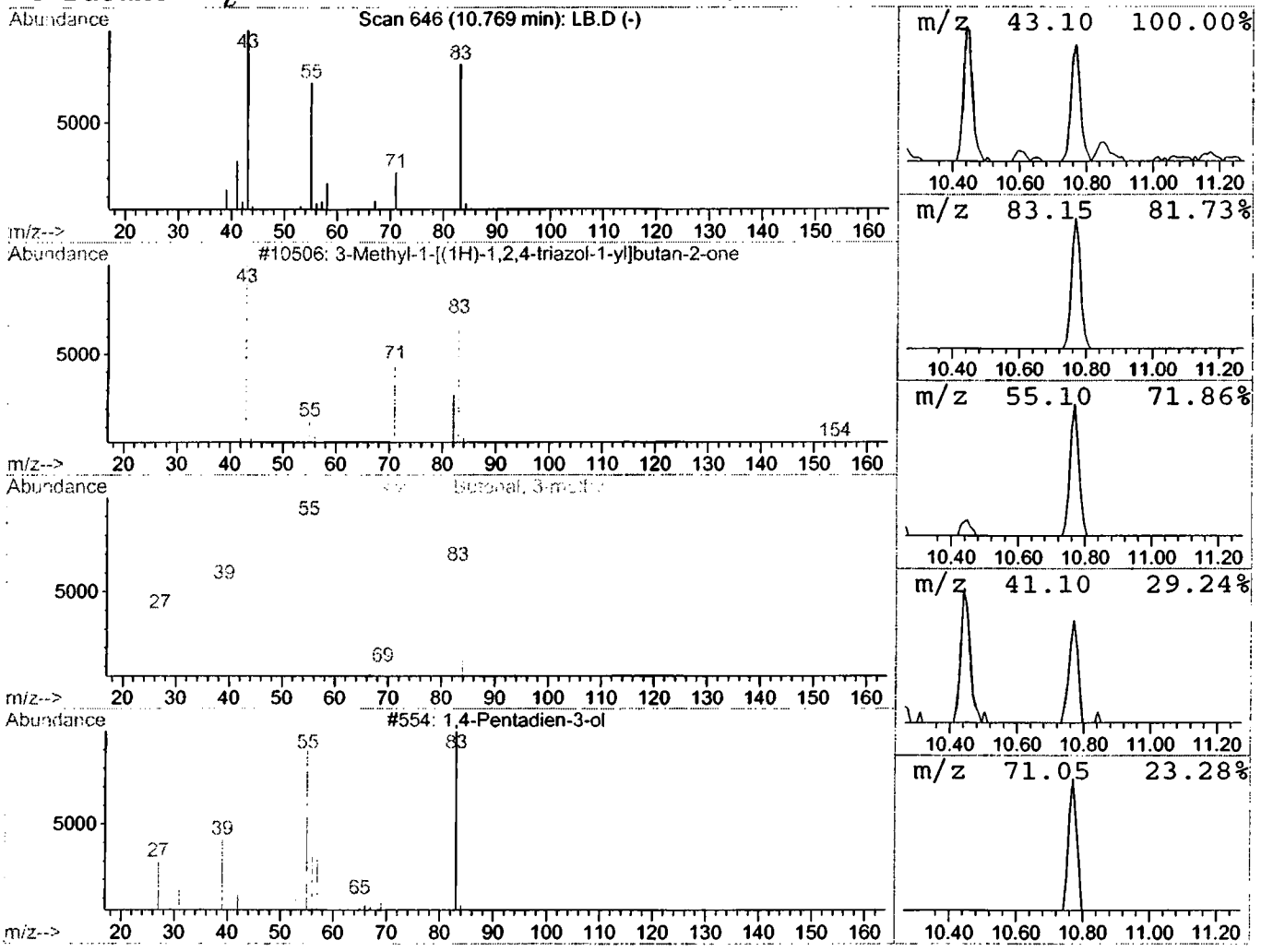
Quant Method : C:\HPCHEM\1\METHODS\NP112701.M (RTE Integrator)
Title : 208 625/TCLP calibration table
Library : C:\DATABASE\NBS75K.L

****

Peak Number 9 3-Methyl-1-[(1H)-1,2,4-triazol Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.77	0.56 ng/uL	120862	1,4-Dichlorobenzene-d4	11.80

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	3-Methyl-1-[(1H)-1,2,4-triazol-1-yl	153	C7H11N3O	064922-02-7	33
2	2-Butenal, 3-methyl-	84	C5H8O	000107-86-8	9
3	1,4-Pentadien-3-ol	84	C5H8O	000922-65-6	9
4	Butane	58	C4H10	000106-97-8	9



Library Search Compound Report

Data File : C:\HPCHEM\1\DATA\DEC2101\LB.D
Acq On : 22 Dec 2001 2:07 am
Sample : gcms unknown 11/23
Misc :
MS Integration Params: LSCINT.P

Vial: 10
Operator: GALOS
Inst : GC/MS Ins
Multiplr: 1.00

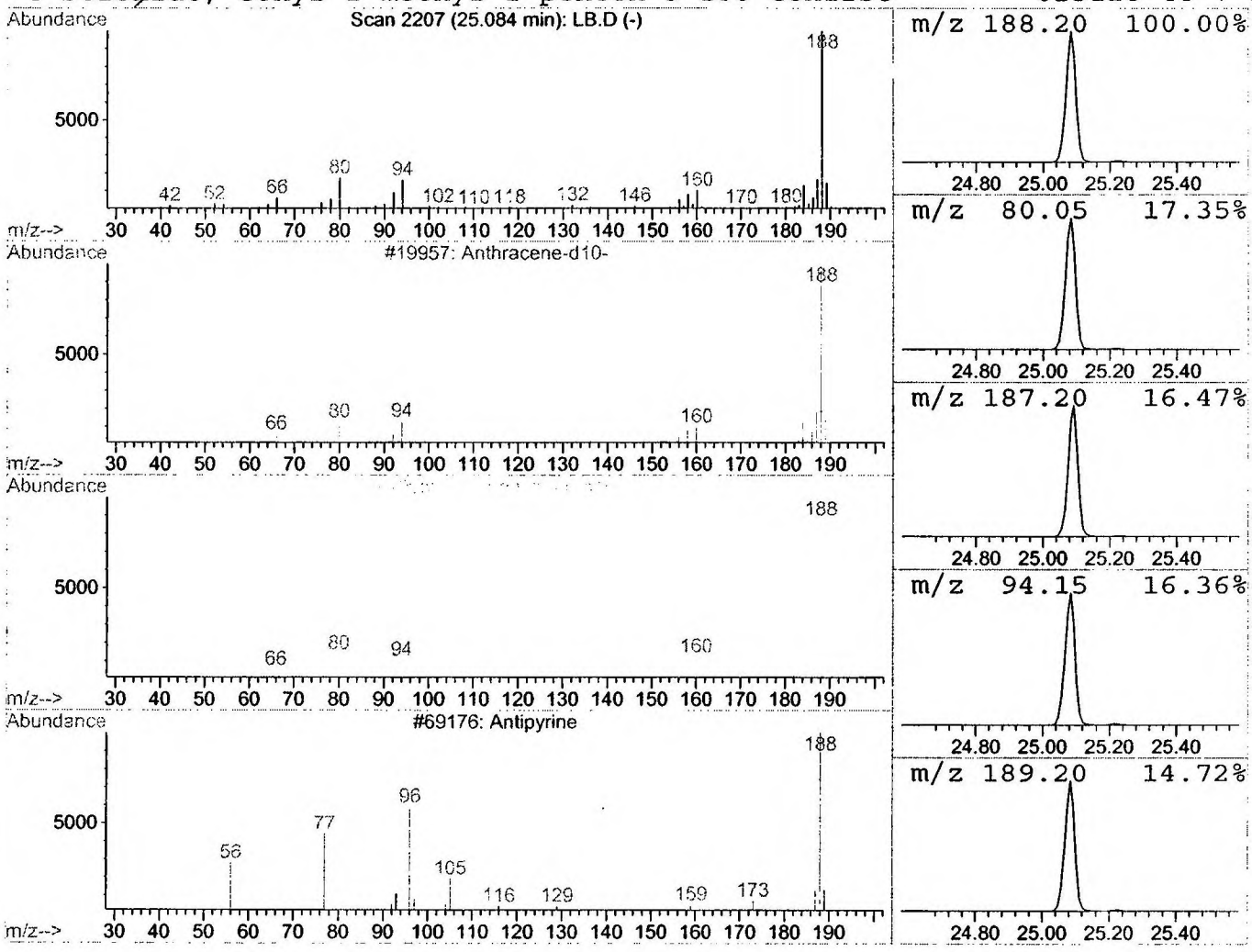
Quant Method : C:\HPCHEM\1\METHODS\NP112701.M (RTE Integrator)
Title : 208 625/TCLP calibration table
Library : C:\DATABASE\NBS75K.L

****

Peak Number 10 Anthracene-d10- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
25.08	20.00 ng/uL	4963960	Phenanthrene-d10	25.22

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Anthracene-d10-	188	C14D10	001719-06-8	96
2	Phenanthrene-d10	188	C14D10	001517-22-2	52
3	Antipyrine	188	C11H12N2O	000060-80-0	36
4	Selenide, ethyl 1-methyl-1-penten-3	188	C8H12Se	025128-48-7	33



Tentatively Identified Compound (LSC) summary

Operator ID: GALOS Date Acquired: 22 Dec 2001 2:07 am
Data File: C:\HPCHEM\1\DATA\DEC2101\LB.D
Name: gcms unknown 11/23
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
1-Propene, 2-fluoro-	5.14	0.5	ng/uL	98014	ISTD01	11.80	4319370	20.0
3-Hexanone	6.40	0.7	ng/uL	144116	ISTD01	11.80	4319370	20.0
2-Hexanone	6.50	1.1	ng/uL	242562	ISTD01	11.80	4319370	20.0
3-Hexanol	6.63	0.5	ng/uL	107868	ISTD01	11.80	4319370	20.0
2-Hexanol	6.74	0.8	ng/uL	163024	ISTD01	11.80	4319370	20.0
2-Pentanone, 4-hydro	7.77	2.1	ng/uL	448535	ISTD01	11.80	4319370	20.0
1-Butanol, 2,2-dimet	10.24	0.4	ng/uL	95992	ISTD01	11.80	4319370	20.0
2-Pentanol, 4-methyl	10.45	0.5	ng/uL	116114	ISTD01	11.80	4319370	20.0
3-Methyl-1-[(1H)-1,2	10.77	0.6	ng/uL	120862	ISTD01	11.80	4319370	20.0
Anthracene-d10-	25.08	20.0	ng/uL	4963960	ISTD04	25.22	4963960	20.0

LB.D NP112701.M Fri Dec 28 11:47:01 2001