Quantitation Report (QT Reviewed)

Data File : C:\HPCHEM\1\DATA\DEC0601\LBB.D
Acq On : 6 Dec 2001 10:00 pm
Sample : lab blank 11/8
Misc :
MS Integration Params: RTEINT.P
Quant Time: Dec 11 11:37 2001

Vial: 14
Operator: GALOS
Inst : GC/MS Ins
Multiplr: 1.00
Quant Results File: NP112701.RES

26833, 26845-7
26982-8
26991-3
27007-9
27011-2

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	1205687	20.00	ng/uL	-0.05
19) Naphthalene-d8	15.50	136	4604078	20.00	ng/uL	-0.05
31) Acenaphthene-d10	20.77	164	2099779	20.00	ng/uL	-0.05
49) Phenanthrene-d10	25.18	188	3055515	20.00	ng/uL	-0.06
59) Chrysene-d12	33.19	240	1992675	20.00	ng/uL	-0.06
68) Perylene-d12	37.28	264	1417144	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	429	0.01	ng/uL	0.46
Spiked Amount 75.000	Range 34 - 84		Recovery = 0.01%#			
7) 1,2-Dichlorobenzene-d4	12.23	152	295	0.01	ng/uL	-0.23
Spiked Amount 50.000	Range 26 - 73		Recovery = 0.02%#			
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	2368	0.02	ng/uL	0.09
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\DEC0601\LBB.D Vial: 14
Acq On : 6 Dec 2001 10:00 pm Operator: GALOS
Sample : lab blank 11/8 Inst : GC/MS Ins
Misc : Multiplr: 1.00
MS Integration Params: RTEINT.P
Quant Time: Dec 11 11:37 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.	
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	22.28	149	330	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	307	N.D.	
56)	Anthracene	25.19	178	307	N.D.	
57)	Di-n-butylphthalate	27.17	149	14085	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\DEC0601\LBB.D Vial: 14
Acq On : 6 Dec 2001 10:00 pm Operator: GALOS
Sample : lab blank 11/8 Inst : GC/MS Ins
Misc : Multiplr: 1.00
MS Integration Params: RTEINT.P
Quant Time: Dec 11 11:37 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	33.19	228	3780	N.D.	
66)	Bis(2-ethylhexyl)phthalate	33.45	149	14851	N.D.	
67)	Chrysene	33.19	228	3780	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	37.29	252	4659	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.	

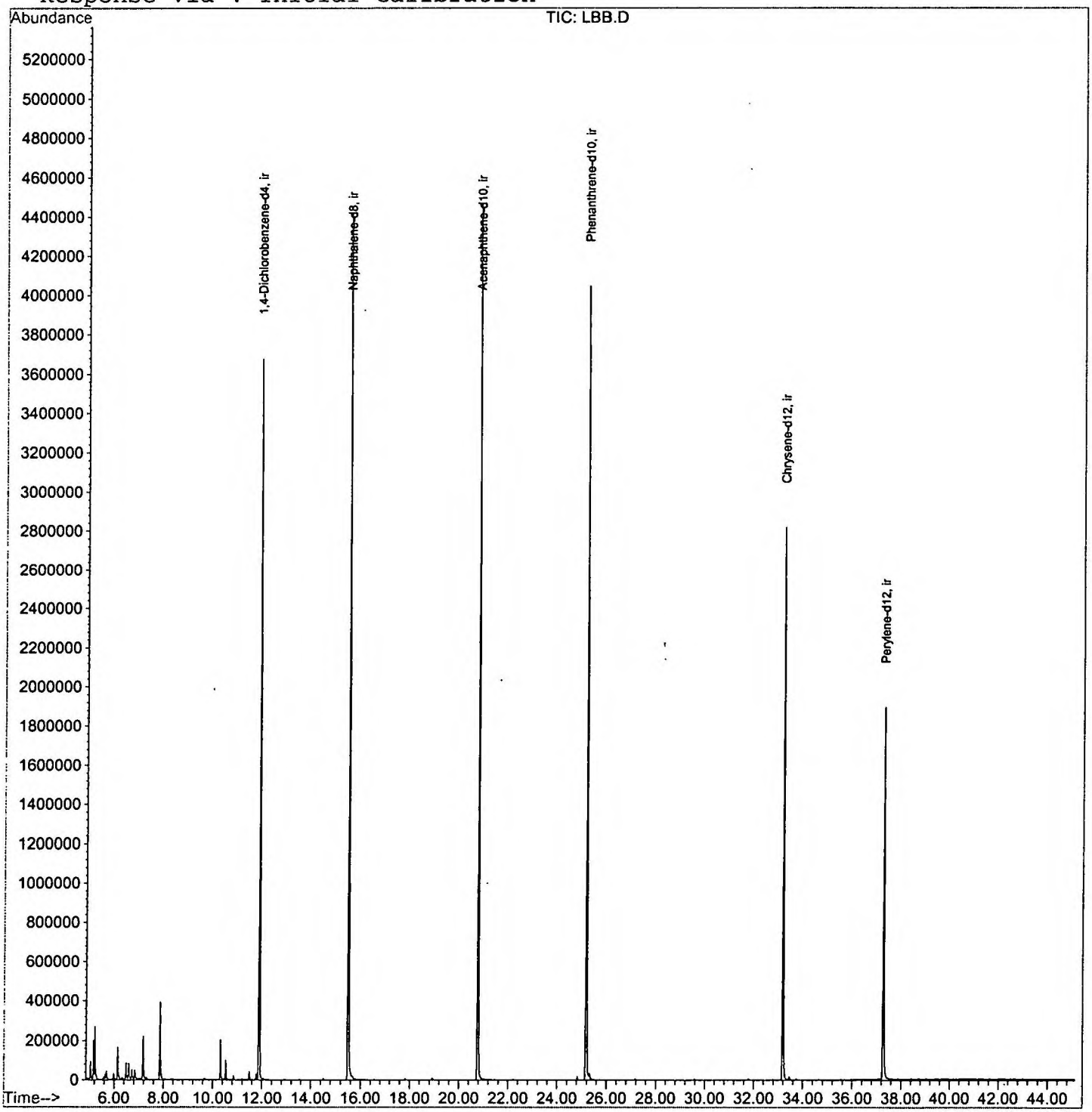
(#) = qualifier out of range (m) = manual integration

Quantitation Report

Data File : C:\HPCHEM\1\DATA\DEC0601\LBB.D
Acq On : 6 Dec 2001 10:00 pm
Sample : lab blank 11/8
Misc :
MS Integration Params: RTEINT.P
Quant Time: Dec 11 11:37 2001

Vial: 14
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\DEC0601\LBB.D
Acq On : 6 Dec 2001 10:00 pm
Sample : lab blank 11/8
Misc :
MS Integration Params: LSCINT.P

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

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

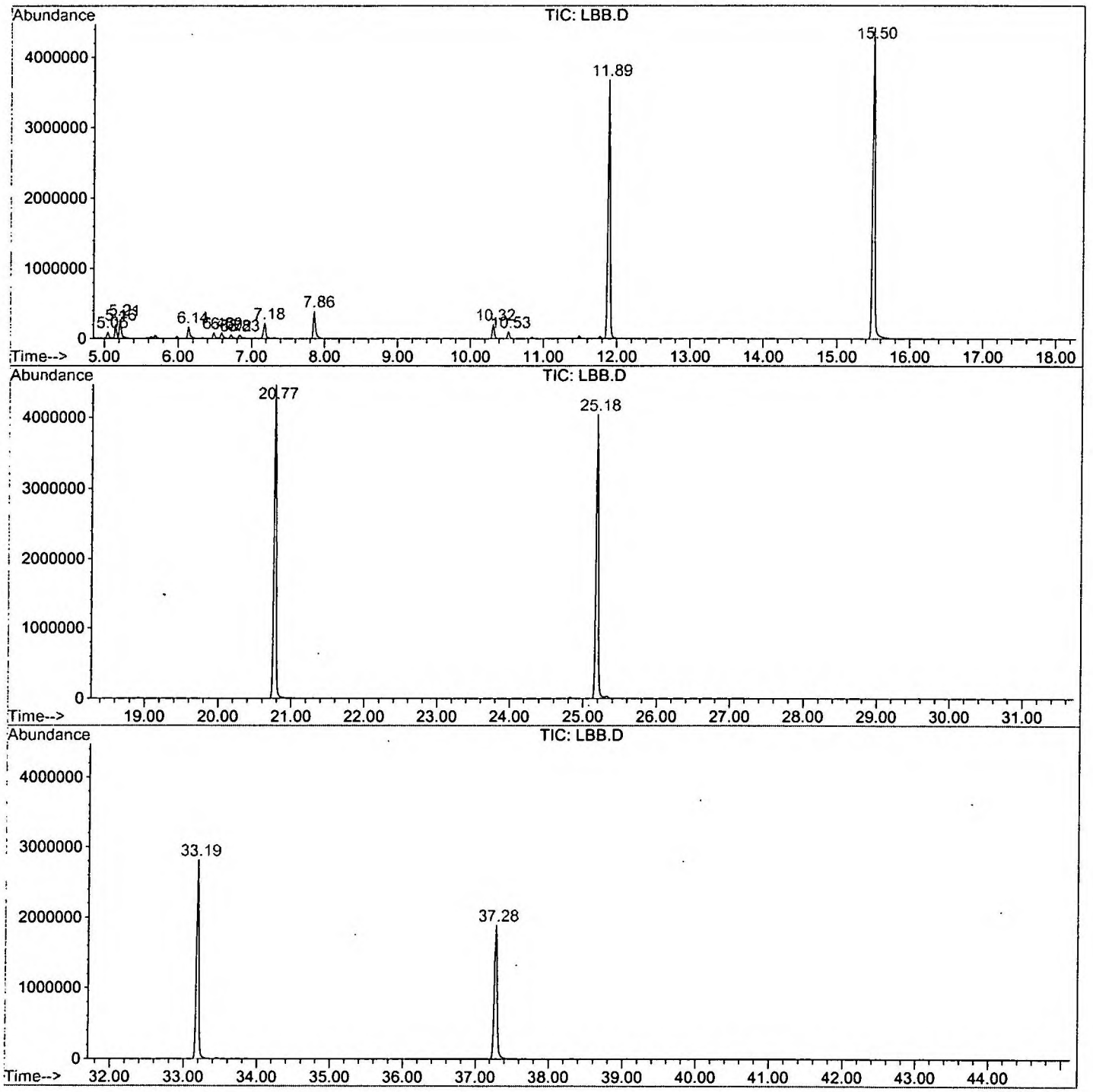
peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	peak area	peak % max.	% of total
1	5.045	15	22	28	rBV	91987	177088	1.90%	0.354%
2	5.155	30	34	37	rVV	198457	330059	3.53%	0.659%
3	5.210	37	40	56	rVB	268307	572320	6.12%	1.143%
4	6.136	137	141	155	rBV	166966	366317	3.92%	0.732%
5	6.485	175	179	187	rBV	86186	170042	1.82%	0.340%
6	6.595	187	191	201	rVV	81189	186218	1.99%	0.372%
7	6.723	201	205	213	rVB	50216	109662	1.17%	0.219%
8	6.833	213	217	228	rBV2	47930	122294	1.31%	0.244%
9	7.182	248	255	263	rBV	220954	504398	5.40%	1.008%
10	7.860	325	329	341	rBV	392100	821835	8.79%	1.642%
11	10.318	592	597	605	rBV	204683	391770	4.19%	0.783%
12	10.529	614	620	630	rBV	100164	199001	2.13%	0.398%
13	11.886	763	768	785	rVB	3674771	7422712	79.43%	14.828%
14	15.499	1155	1162	1181	rBV	4420310	9344433	100.00%	18.667%
15	20.772	1731	1737	1754	rBV	4470753	9231935	98.80%	18.442%
16	25.183	2212	2218	2230	rBV2	4045922	8568909	91.70%	17.118%
17	33.188	3084	3091	3108	rBV2	2817559	6353289	67.99%	12.692%
18	37.278	3529	3537	3554	rBV2	1889581	5187027	55.51%	10.362%

Sum of corrected areas: 50059309

LBB.D NP112701.M Wed Dec 12 08:18:53 2001

LSC Report - Integrated Chromatogram

File : C:\HPCHEM\1\DATA\DEC0601\LBB.D
Operator : GALOS
Acquired : 6 Dec 2001 10:00 pm using AcqMethod NP112701
Instrument : GC/MS Ins
Sample Name: lab blank 11/8
Misc Info :
Vial Number: 14
Quant File :NP112701.RES (RTE Integrator)



LBB.D NP112701.M Wed Dec 12 08:18:56 2001

Library Search Compound Report

Data File : C:\HPCHEM\1\DATA\DEC0601\LBB.D
Acq On : 6 Dec 2001 10:00 pm
Sample : lab blank 11/8
Misc :
MS Integration Params: LSCINT.P

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

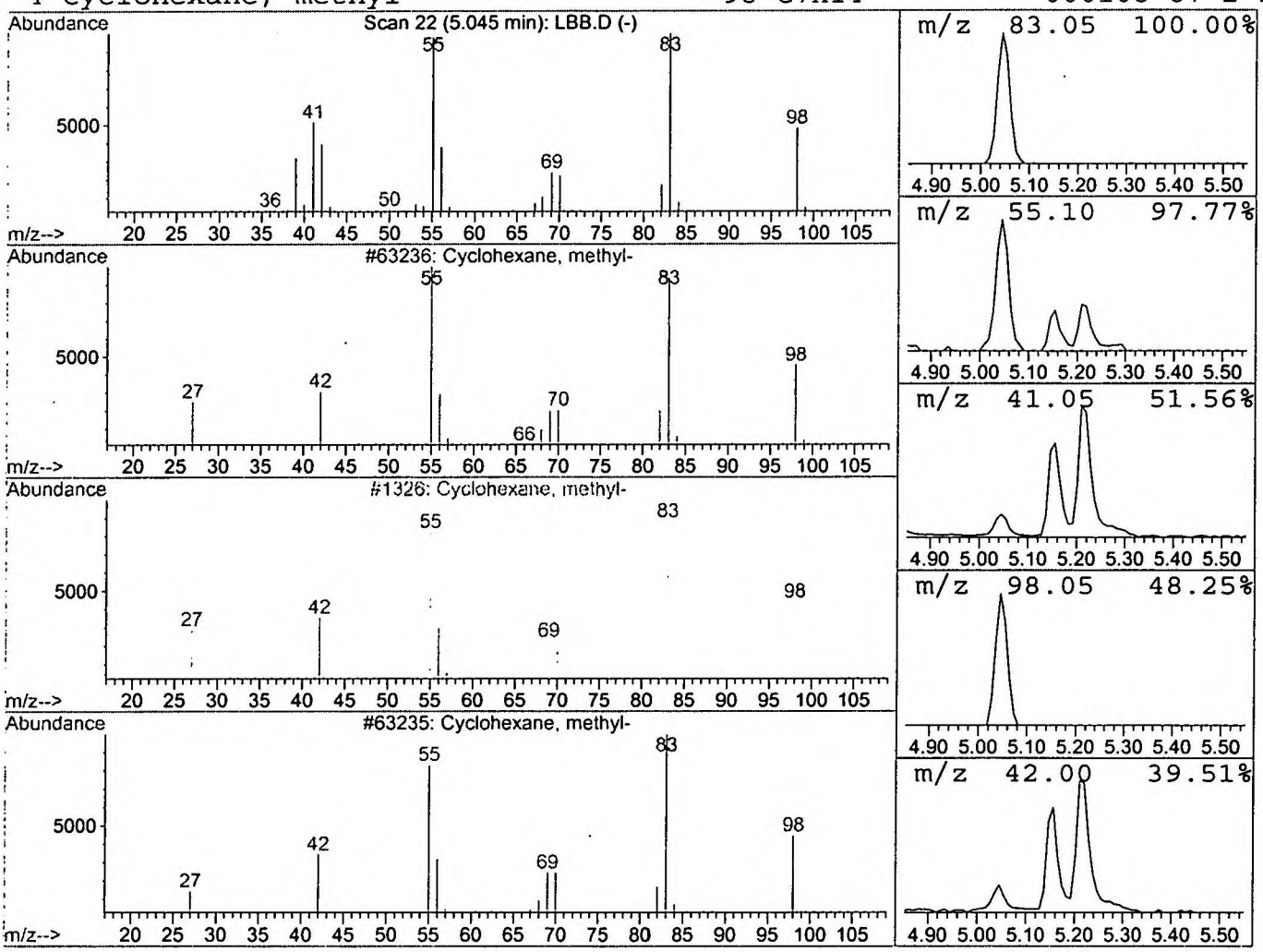
Quant Method : C:\HPCHEM\1\METHODS\NP112701.M (RTE Integrator)
Title : 208 625/TCLP calibration table
Library : C:\DATABASE\NBS75K.L

****

Peak Number 1 Cyclohexane, methyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.05	0.48 ng/uL	177088	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	97
3	Cyclohexane, methyl-	98	C7H14	000108-87-2	97
4	Cyclohexane, methyl-	98	C7H14	000108-87-2	53



Library Search Compound Report

Data File : C:\HPCHEM\1\DATA\DEC0601\LBB.D
Acq On : 6 Dec 2001 10:00 pm
Sample : lab blank 11/8
Misc :
MS Integration Params: LSCINT.P

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

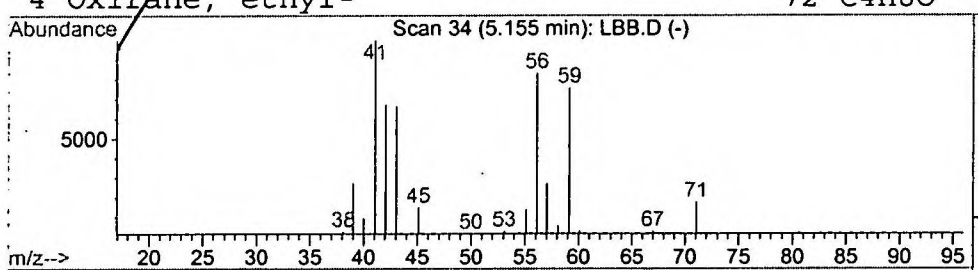
Quant Method : C:\HPCHEM\1\METHODS\NP112701.M (RTE Integrator)
Title : 208 625/TCLP calibration table
Library : C:\DATABASE\NBS75K.L

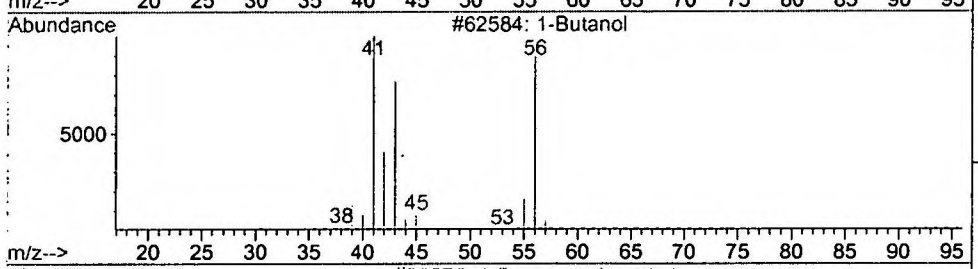
****

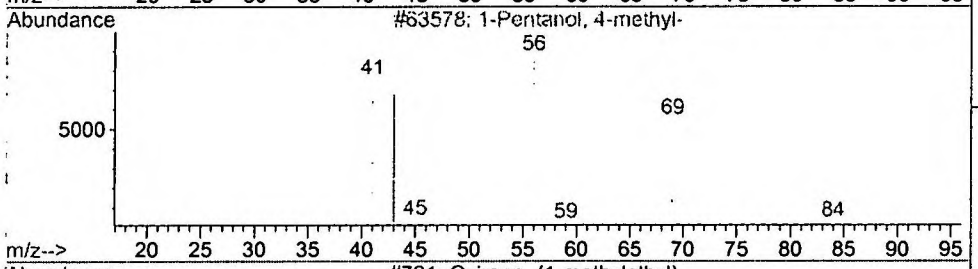
Peak Number 2 1-Butanol Concentration Rank 6

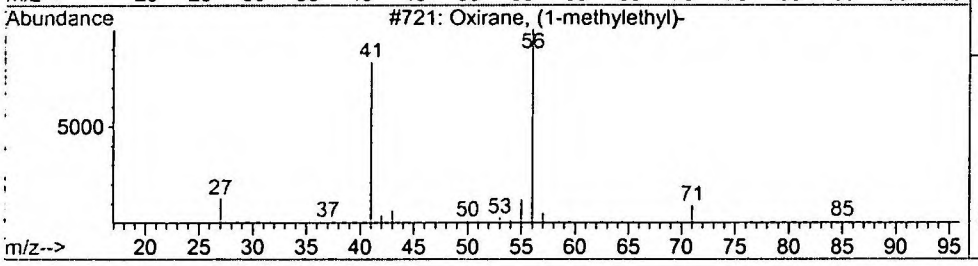
R.T.	EstConc	Area	Relative to ISTD	R.T.
5.16	0.89 ng/uL	330059	1,4-Dichlorobenzene-d4	11.89

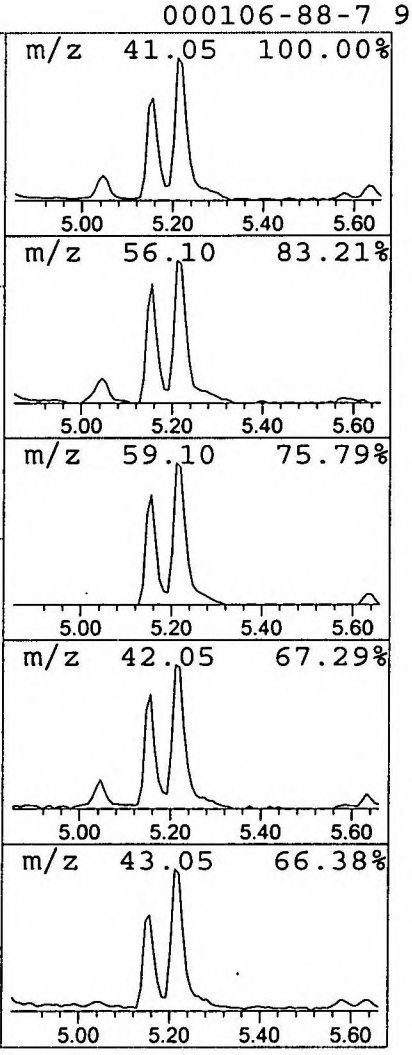
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Butanol	74	C4H10O	000071-36-3	32
2	1-Pentanol, 4-methyl-	102	C6H14O	000626-89-1	25
3	Oxirane, (1-methylethyl)-	86	C5H10O	001438-14-8	17
4	Oxirane, ethyl-	72	C4H8O	000106-88-7	9











Library Search Compound Report

Data File : C:\HPCHEM\1\DATA\DEC0601\LBB.D
Acq On : 6 Dec 2001 10:00 pm
Sample : lab blank 11/8
Misc :
MS Integration Params: LSCINT.P

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

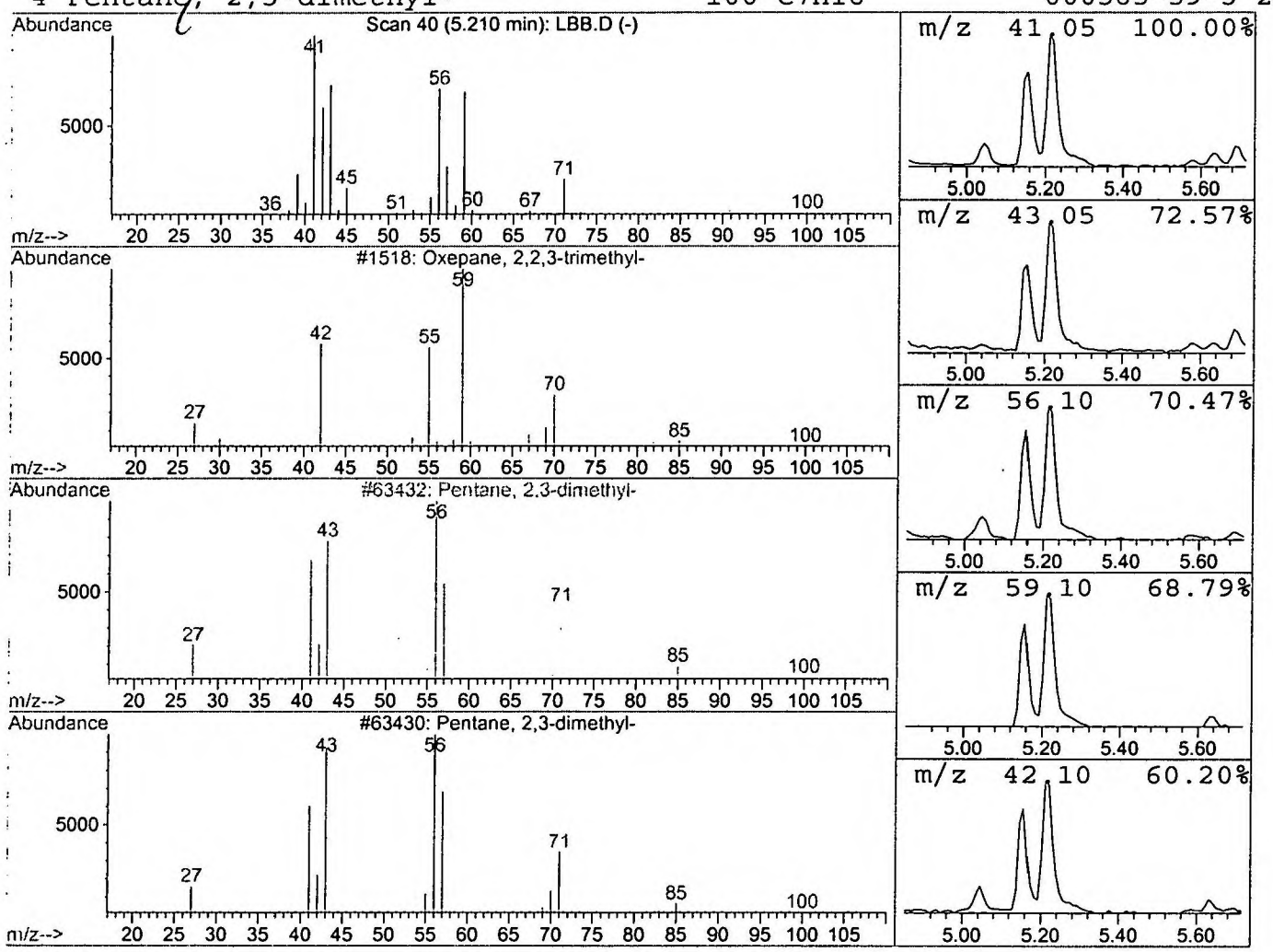
Quant Method : C:\HPCHEM\1\METHODS\NP112701.M (RTE Integrator)
Title : 208 625/TCLP calibration table
Library : C:\DATABASE\NBS75K.L

****

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.21	1.54 ng/uL	572320	1,4-Dichlorobenzene-d4	11.89

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Oxepane, 2,2,3-trimethyl-	100	C6H12O	023120-43-6	35
2	Pentane, 2,3-dimethyl-	100	C7H16	000565-59-3	35
3	Pentane, 2,3-dimethyl-	100	C7H16	000565-59-3	35
4	Pentane, 2,3-dimethyl-	100	C7H16	000565-59-3	25



Library Search Compound Report

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Acq On : 6 Dec 2001 10:00 pm
Sample : lab blank 11/8
Misc :
MS Integration Params: LSCINT.P

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

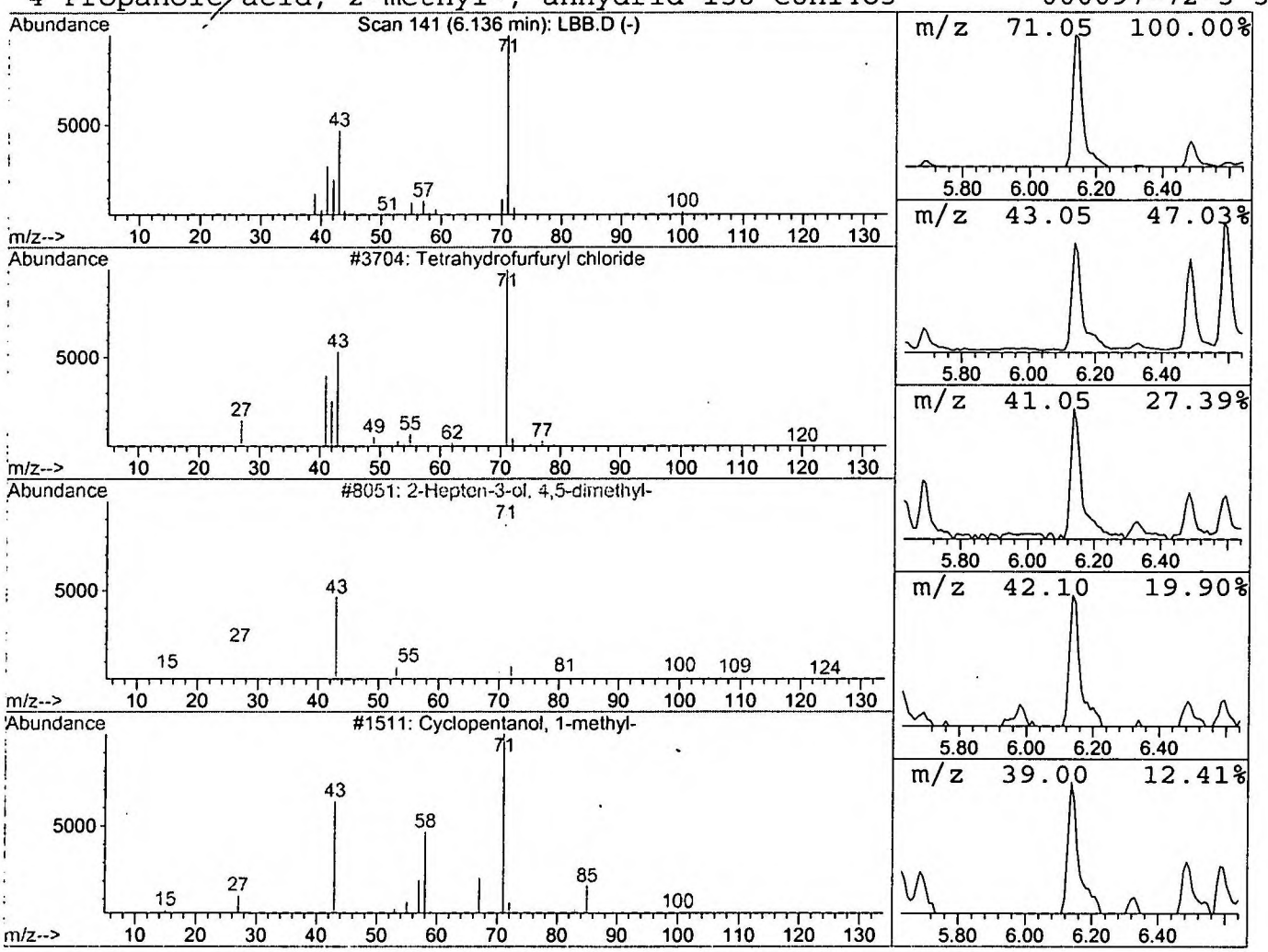
Quant Method : C:\HPCHEM\1\METHODS\NP112701.M (RTE Integrator)
Title : 208 625/TCLP calibration table
Library : C:\DATABASE\NBS75K.L

****

Peak Number 4 Tetrahydrofurfuryl chloride Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.14	0.99 ng/uL	366317	1,4-Dichlorobenzene-d4	11.89

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Tetrahydrofurfuryl chloride	120	C5H9ClO	003003-84-7	50
2	2-Hepten-3-ol, 4,5-dimethyl-	142	C9H18O	055956-37-1	39
3	Cyclopentanol, 1-methyl-	100	C6H12O	001462-03-9	36
4	Propanoic acid, 2-methyl-, anhydrid	158	C8H14O3	000097-72-3	33



Library Search Compound Report

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Acq On : 6 Dec 2001 10:00 pm
Sample : lab blank 11/8
Misc :
MS Integration Params: LSCINT.P

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

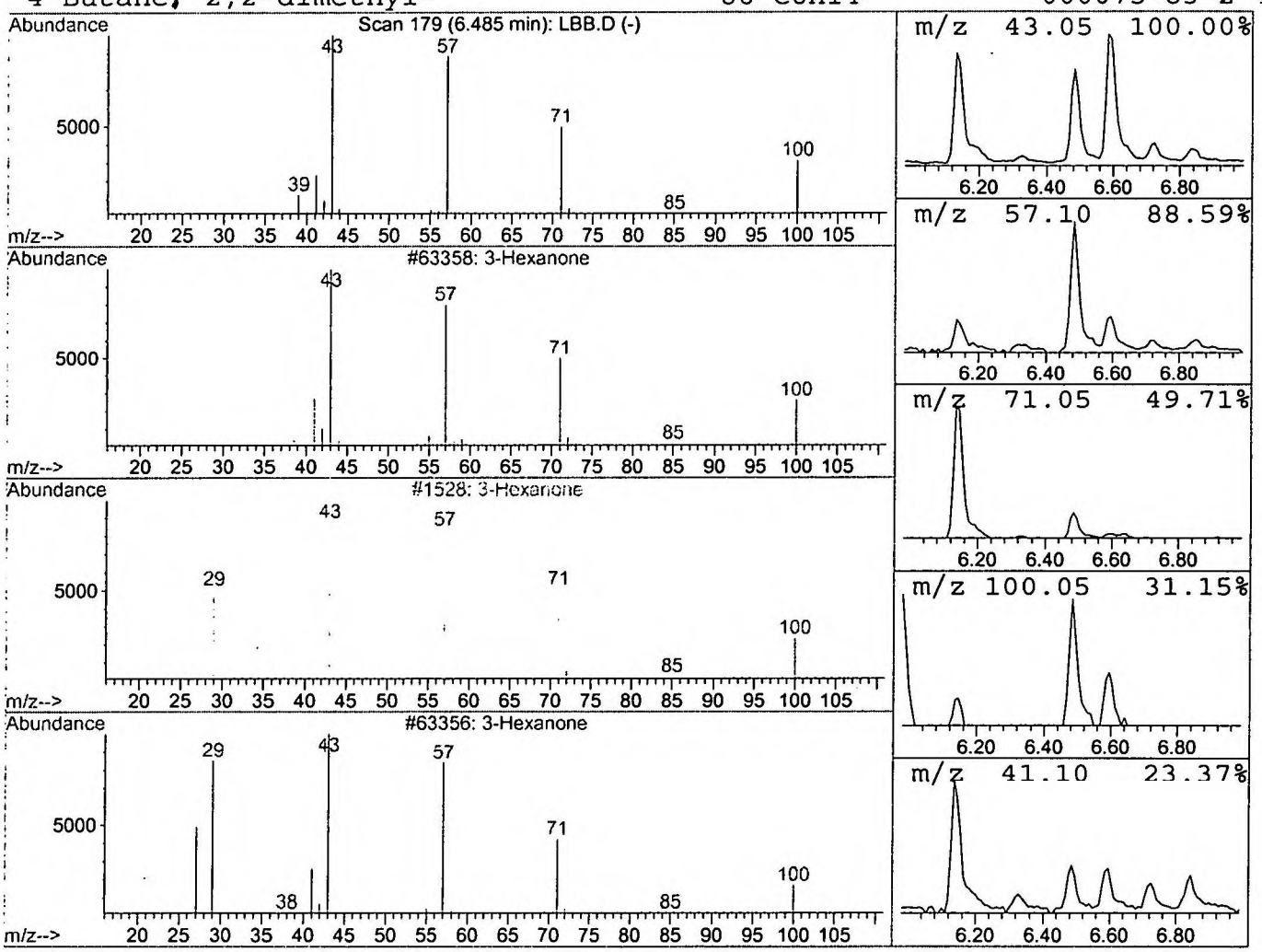
Quant Method : C:\HPCHEM\1\METHODS\NP112701.M (RTE Integrator)
Title : 208 625/TCLP calibration table
Library : C:\DATABASE\NBS75K.L

****

Peak Number 5 3-Hexanone Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.48	0.46 ng/uL	170042	1,4-Dichlorobenzene-d4	11.89

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	91
3	3-Hexanone	100	C6H12O	000589-38-8	78
4	Butane, 2,2-dimethyl-	86	C6H14	000075-83-2	42



Library Search Compound Report

Data File : C:\HPCHEM\1\DATA\DEC0601\LBB.D
Acq On : 6 Dec 2001 10:00 pm
Sample : lab blank 11/8
Misc :
MS Integration Params: LSCINT.P

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

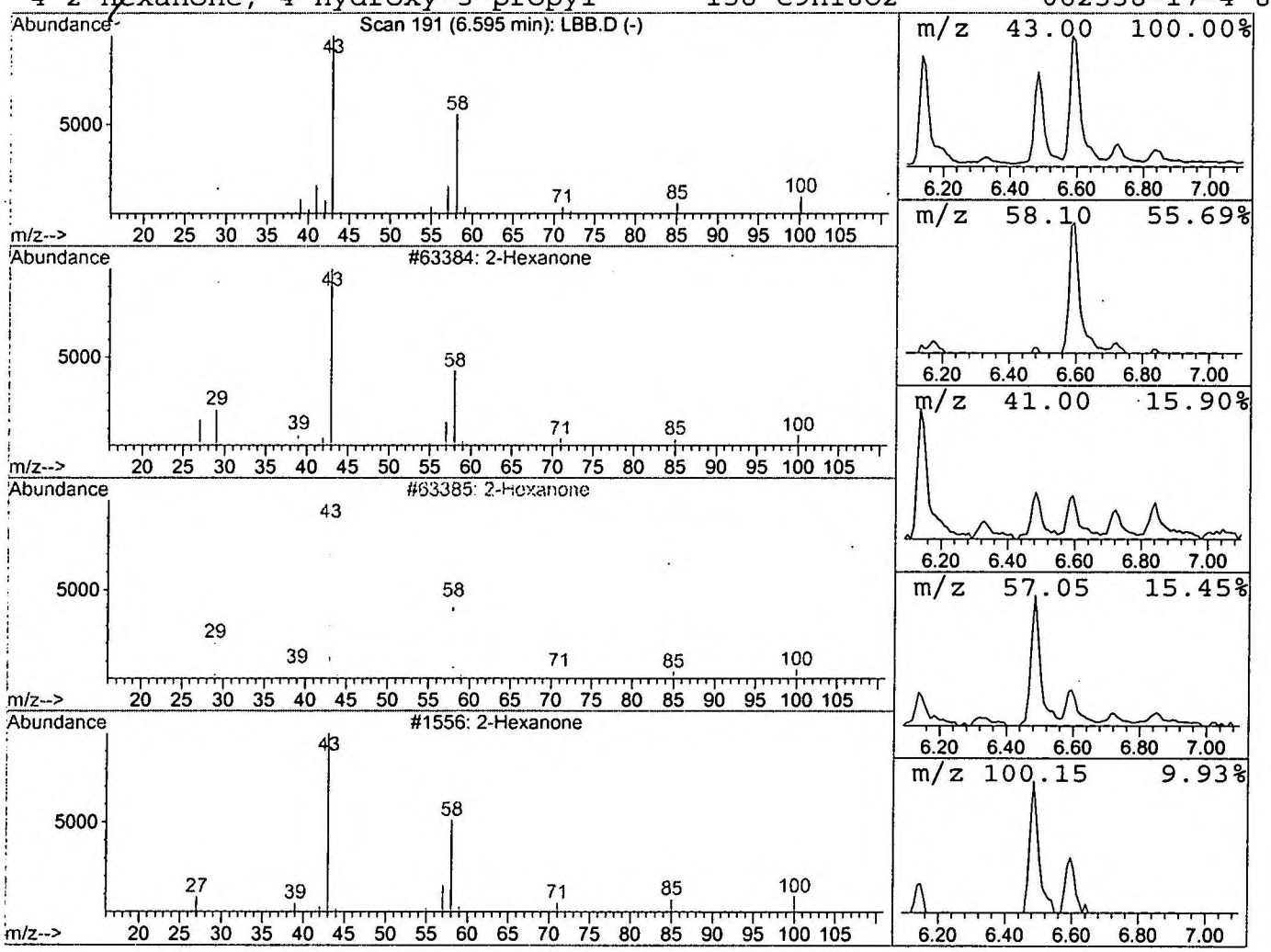
Quant Method : C:\HPCHEM\1\METHODS\NP112701.M (RTE Integrator)
Title : 208 625/TCLP calibration table
Library : C:\DATABASE\NBS75K.L

****

Peak Number 6 2-Hexanone — Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.59	0.50 ng/uL	186218	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	90
3	2-Hexanone	100	C6H12O	000591-78-6	90
4	2-Hexanone, 4-hydroxy-3-propyl-	158	C9H18O2	062338-17-4	83



Library Search Compound Report

Data File : C:\HPCHEM\1\DATA\DEC0601\LBB.D
Acq On : 6 Dec 2001 10:00 pm
Sample : lab blank 11/8
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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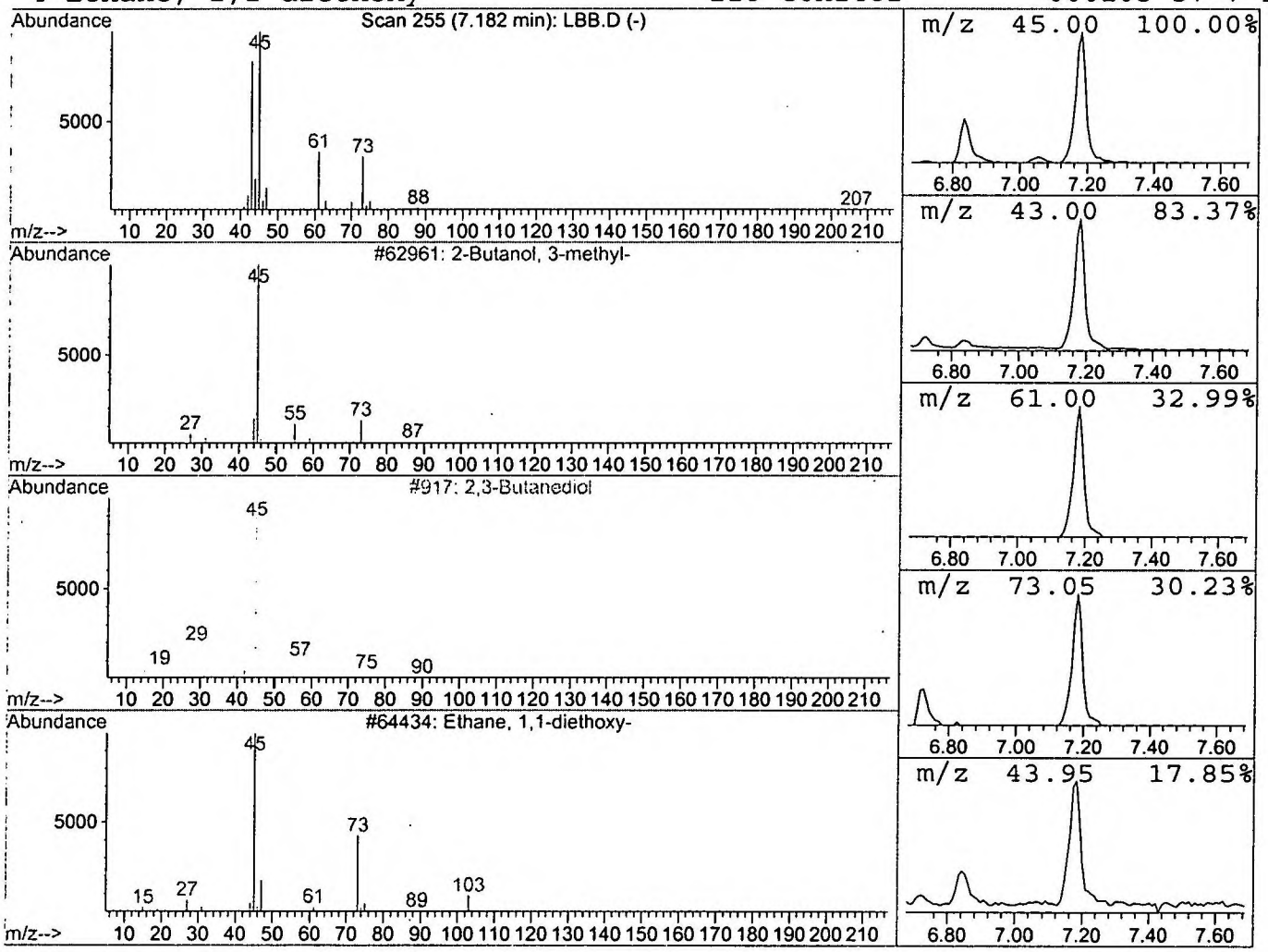
Peak Number 7 2-Butanol, 3-methyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.18	1.36 ng/uL	504398	1,4-Dichlorobenzene-d4	11.89

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Butanol, 3-methyl-	88	C5H12O	000598-75-4	38
2	2,3-Butanediol	90	C4H10O2	000513-85-9	28
3	Ethane, 1,1-diethoxy-	118	C6H14O2	000105-57-7	9
4	Ethane, 1,1-diethoxy-	118	C6H14O2	000105-57-7	9



Library Search Compound Report

Data File : C:\HPCHEM\1\DATA\DEC0601\LBB.D Vial: 14
Acq On : 6 Dec 2001 10:00 pm Operator: GALOS
Sample : lab blank 11/8 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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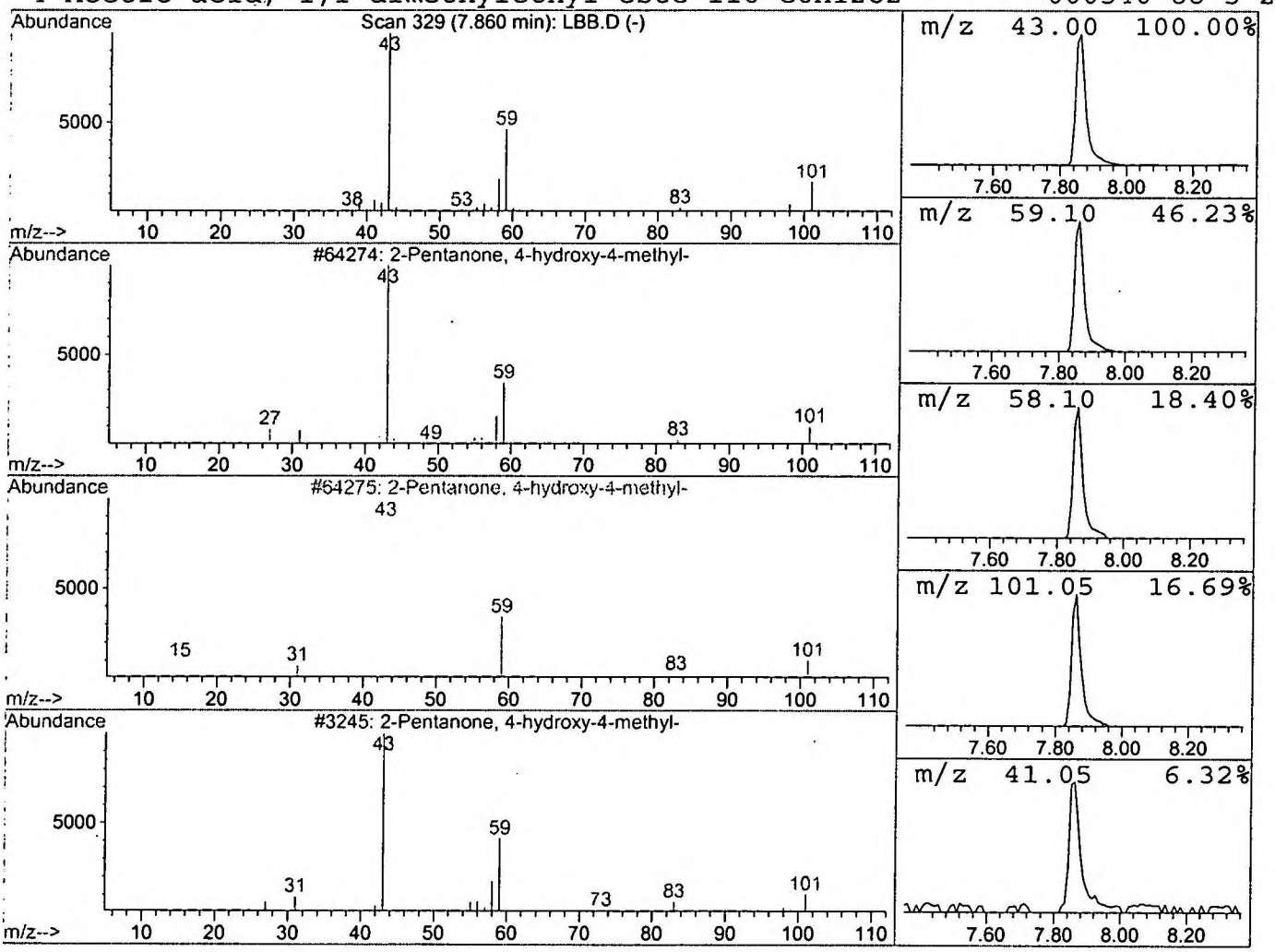
Peak Number 8 2-Pentanone, 4-hydroxy-4-methy Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.86	2.21 ng/uL	821835	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	83
2	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	59
3	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	43
4	Acetic acid, 1,1-dimethylethyl este	116	C6H12O2	000540-88-5	28



Library Search Compound Report

Data File : C:\HPCHEM\1\DATA\DEC0601\LBB.D
Acq On : 6 Dec 2001 10:00 pm
Sample : lab blank 11/8
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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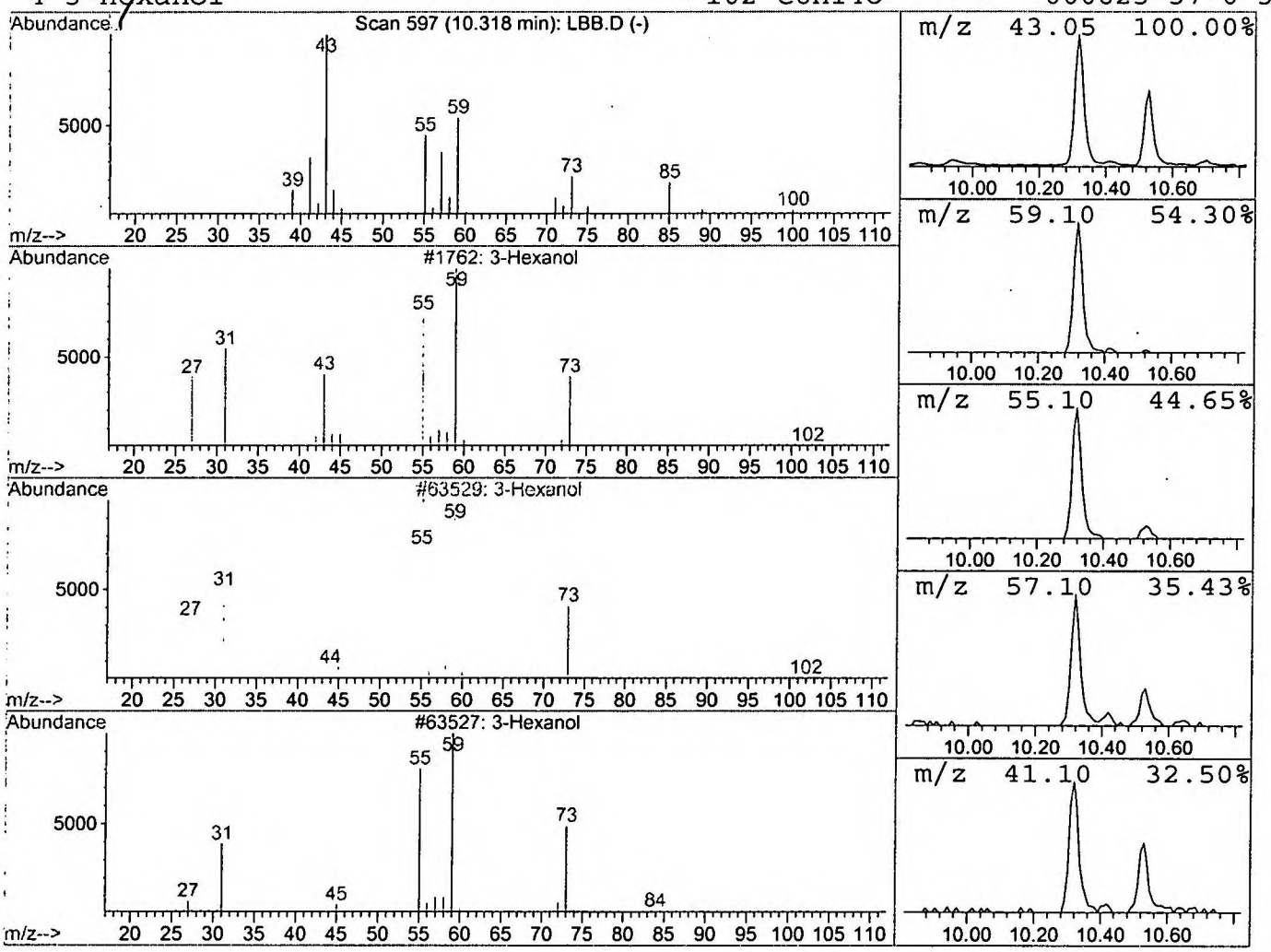
Peak Number 9 3-Hexanol Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.32	1.06 ng/uL	391770	1,4-Dichlorobenzene-d4	11.89

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	3-Hexanol	102	C6H14O	000623-37-0	38
2	3-Hexanol	102	C6H14O	000623-37-0	37
3	3-Hexanol	102	C6H14O	000623-37-0	37
4	3-Hexanol	102	C6H14O	000623-37-0	32



Library Search Compound Report

Data File : C:\HPCHEM\1\DATA\DEC0601\LBB.D
Acq On : 6 Dec 2001 10:00 pm
Sample : lab blank 11/8
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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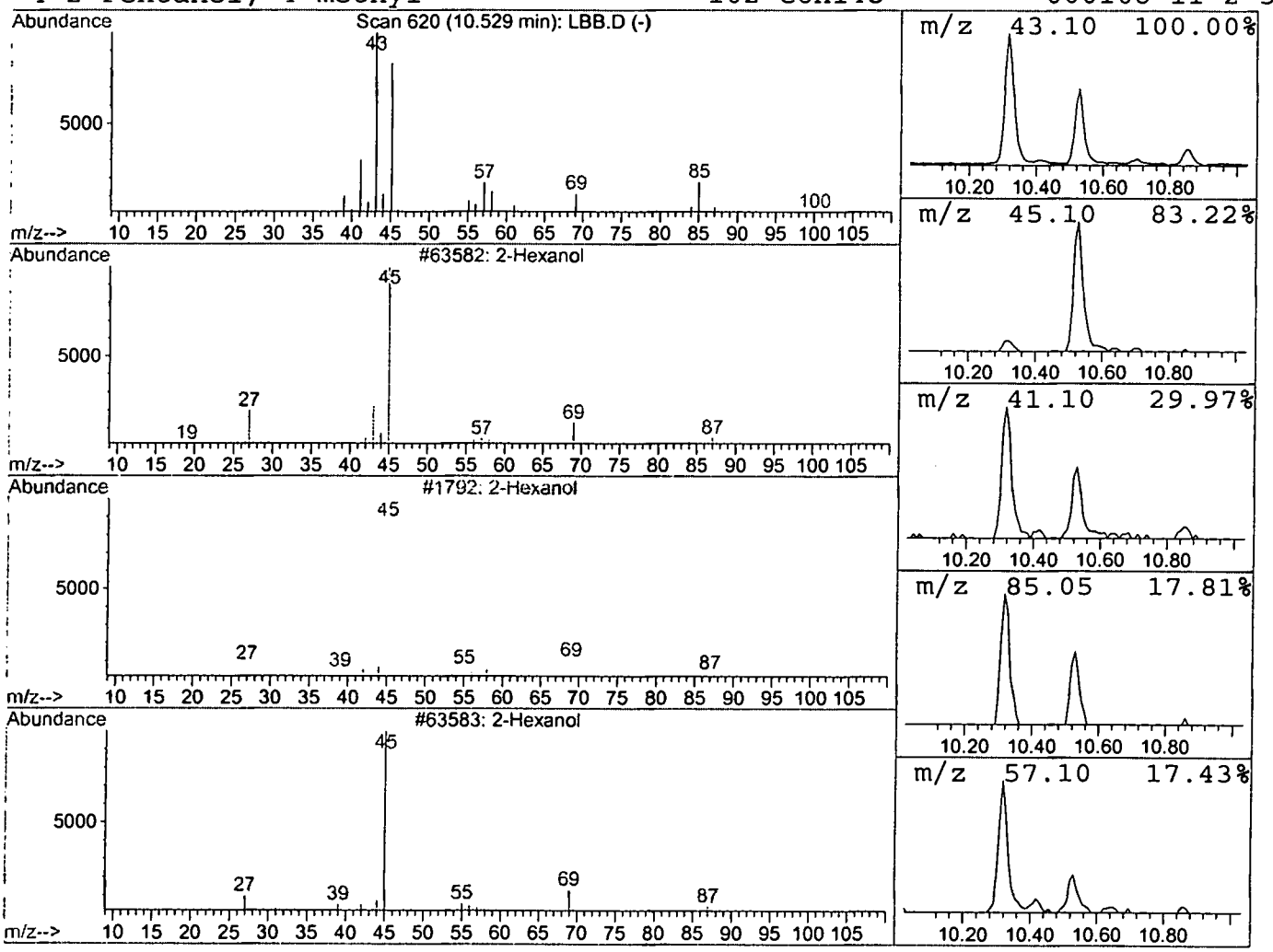
Peak Number 10 2-Hexanol Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.53	0.54 ng/uL	199001	1,4-Dichlorobenzene-d4	11.89

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Hexanol	102	C6H14O	000626-93-7	64
2	2-Hexanol	102	C6H14O	000626-93-7	59
3	2-Hexanol	102	C6H14O	000626-93-7	50
4	2-Pentanol, 4-methyl-	102	C6H14O	000108-11-2	50



Tentatively Identified Compound (LSC) summary

Operator ID: GALOS Date Acquired: 6 Dec 2001 10:00 pm
Data File: C:\HPCHEM\1\DATA\DEC0601\LBB.D
Name: lab blank 11/8
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-	5.05	0.5	ng/uL	177088	ISTD01	11.89	7422710	20.0
1-Butanol	5.16	0.9	ng/uL	330059	ISTD01	11.89	7422710	20.0
Oxepane, 2,2,3-trime	5.21	1.5	ng/uL	572320	ISTD01	11.89	7422710	20.0
Tetrahydrofurfuryl c	6.14	1.0	ng/uL	366317	ISTD01	11.89	7422710	20.0
3-Hexanone	6.48	0.5	ng/uL	170042	ISTD01	11.89	7422710	20.0
2-Hexanone	6.59	0.5	ng/uL	186218	ISTD01	11.89	7422710	20.0
2-Butanol, 3-methyl-	7.18	1.4	ng/uL	504398	ISTD01	11.89	7422710	20.0
2-Pentanone, 4-hydro	7.86	2.2	ng/uL	821835	ISTD01	11.89	7422710	20.0
3-Hexanol	10.32	1.1	ng/uL	391770	ISTD01	11.89	7422710	20.0
2-Hexanol	10.53	0.5	ng/uL	199001	ISTD01	11.89	7422710	20.0

LBB.D NP112701.M Wed Dec 12 08:19:18 2001