Quantitation Report (QT Reviewed)

Data File : C:\HPCHEM\1\DATA\OCT0201\LB.D
Acq On : 2 Oct 2001 3:06 pm
Sample : 9/20
Misc :
MS Integration Params: RTEINT.P
Quant Time: Oct 3 9:13 2001

Lab Blank 22779, 22913-22919 22974-22978 22995-22998

Vial: 2
Operator: 209 GALOS
Inst : GC/MS 209
Multiplr: 1.00
Quant Results File: NP091101.RES

Quant Method : C:\HPCHEM\1\METHODS\NP091101.M (RTE Integrator)
Title : 209 625/TCLP calibration table 7/16/01
Last Update : Thu Sep 13 12:47:04 2001
Response via : Initial Calibration
DataAcq Meth : NP091101

See last Page

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	11.78	152	496271	20.00	ug/l	-0.12
19) Naphthalene-d8	15.40	136	1958686	20.00	ug/l	-0.12
31) Acenaphthene-d10	20.69	164	931392	20.00	ug/l	-0.12
49) Phenanthrene-d10	25.11	188	1332063	20.00	ug/l	-0.12
59) Chrysene-d12	33.16	240	980303	20.00	ug/l	-0.12
68) Perylene-d12	37.28	264	747027	20.00	ug/l	-0.14
System Monitoring Compounds						
4) 2-Fluorophenol	0.00	112	0	0.00	ug/l	
Spiked Amount 75.000			Recovery =	0.00%		
5) Phenol-d5	0.00	99	0	0.00	ug/l	
Spiked Amount 75.000			Recovery =	0.00%		
6) 2-Chlorophenol-d4	11.78	132	244	0.01	ug/l	0.37
Spiked Amount 75.000			Recovery =	0.01%		
7) 1,2-Dichlorobenzene-d4	12.30	152	5189	0.21	ug/l	-0.13
Spiked Amount 50.000			Recovery =	0.42%		
20) Nitrobenzene-d5	0.00	82	0	0.00	ug/l	
Spiked Amount 50.000			Recovery =	0.00%		
32) 2-Fluorobiphenyl	18.83	172	1943	0.03	ug/l	0.01
Spiked Amount 50.000			Recovery =	0.06%		
33) 2,4,6-Tribromophenol	0.00	330	0	0.00	ug/l	
Spiked Amount 75.000			Recovery =	0.00%		
62) Terphenyl-d14	0.00	244	0	0.00	ug/l	
Spiked Amount 50.000			Recovery =	0.00%		
Target Compounds						Qvalue
2) Pyridine	5.32	79	281	N.D.		
3) N-nitroso-dimethyl amine	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.		
17) N-Nitroso-di-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

Quantitation Report

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Sample : 9/20
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Title : 209 625/TCLP calibration table 7/16/01
Last Update : Thu Sep 13 12:47:04 2001
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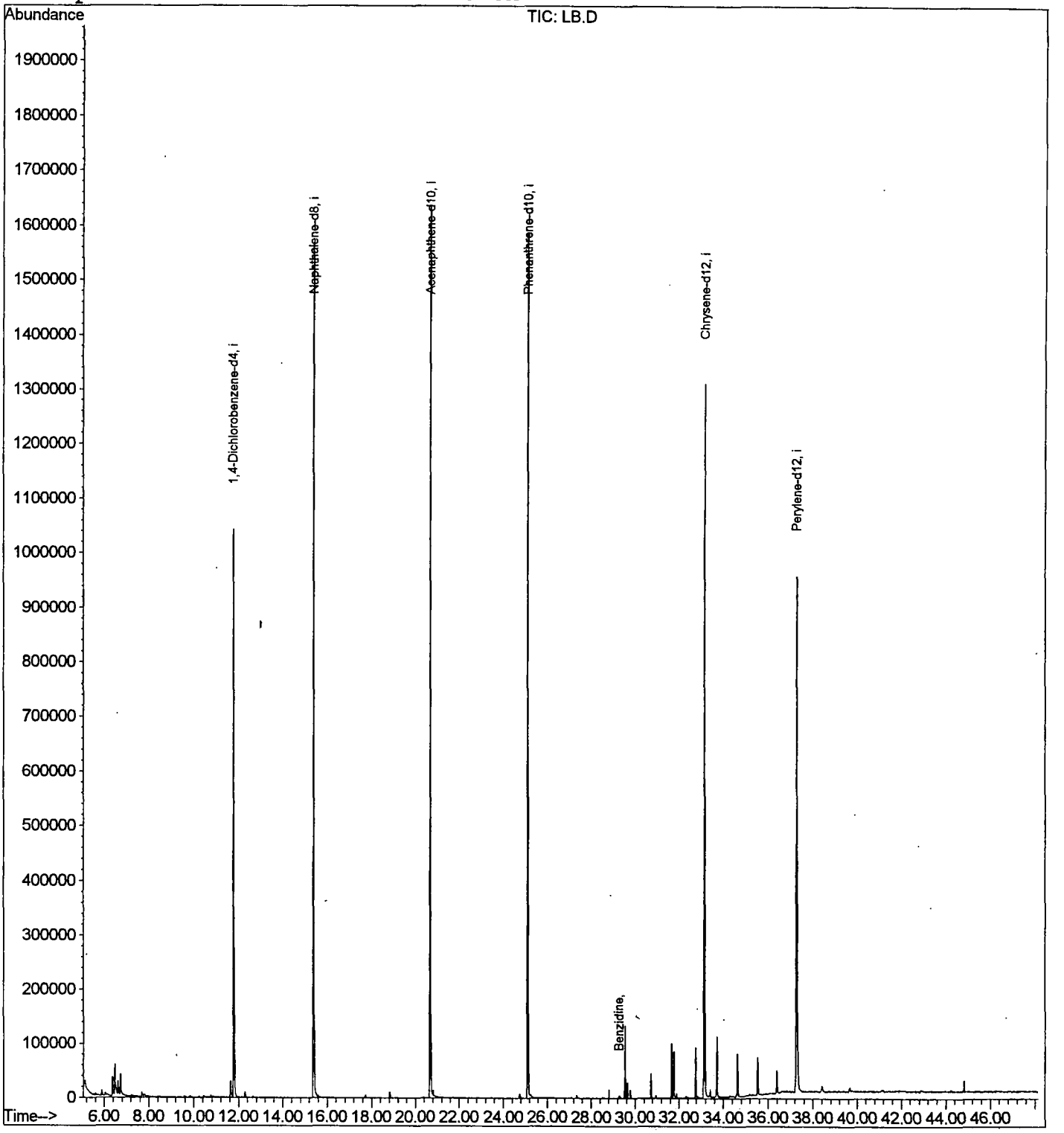


LSC Area Percent Report

Data File : C:\HPCHEM\1\DATA\OCT0201\LB.D
Acq On : 2 Oct 2001 3:06 pm
Sample : 9/20
Misc :
MS Integration Params: LSCINT.P

Vial: 2
Operator: 209 GALOS
Inst : GC/MS 209
Multiplr: 1.00

Method : C:\HPCHEM\1\METHODS\NP091101.M (RTE Integrator)
Title : 209 625/TCLP calibration table 7/16/01
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

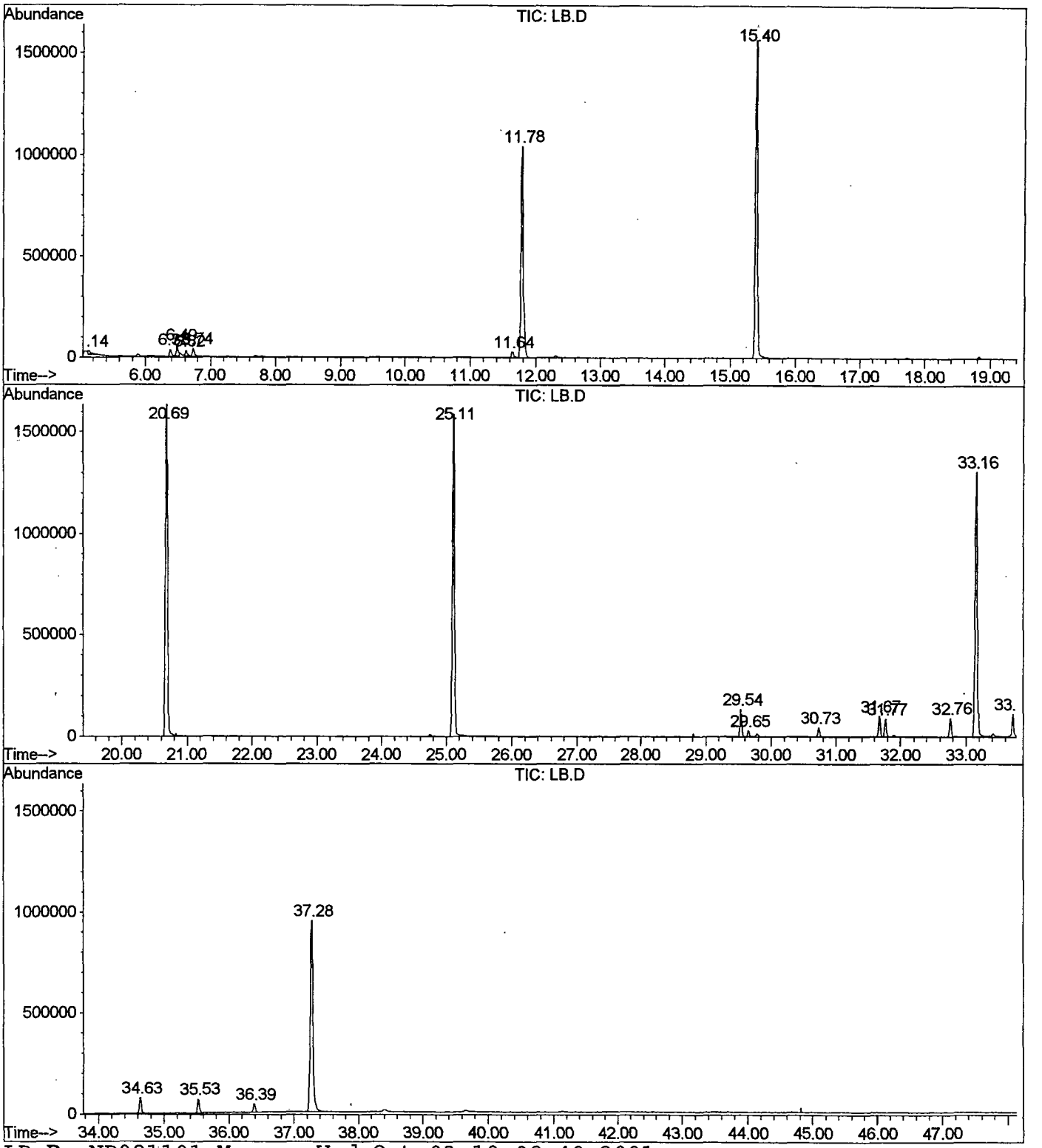
peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	peak area	peak % max.	% of total
1	5.138	9	11	23	rVB	17970	46970	1.27%	0.222%
2	6.378	142	146	153	rBV	34333	76825	2.08%	0.363%
3	6.488	153	158	161	rBV	54801	107024	2.90%	0.506%
4	6.616	169	172	181	rVB	24613	53199	1.44%	0.252%
5	6.736	181	185	198	rVB	38000	88381	2.39%	0.418%
6	11.638	715	719	729	rVB	28778	69043	1.87%	0.326%
7	11.785	729	735	750	rBV	1042087	2566396	69.49%	12.136%
8	15.402	1122	1129	1146	rBV	1565980	3567297	96.59%	16.869%
9	20.690	1698	1705	1718	rBV	1635893	3693277	100.00%	17.465%
10	25.115	2180	2187	2199	rBV2	1586051	3443890	93.25%	16.286%
11	29.539	2663	2669	2675	rBV	133513	260921	7.06%	1.234%
12	29.650	2675	2681	2686	rBV	28538	54299	1.47%	0.257%
13	30.733	2793	2799	2805	rVB	44804	84270	2.28%	0.399%
14	31.669	2893	2901	2907	rBV	100989	204617	5.54%	0.968%
15	31.770	2907	2912	2922	rVB	85552	179169	4.85%	0.847%
16	32.762	3014	3020	3025	rBV	92217	183463	4.97%	0.868%
17	33.157	3056	3063	3078	rBV2	1308298	3048404	82.54%	14.415%
18	33.717	3117	3124	3131	rBV	110571	221588	6.00%	1.048%
19	34.635	3217	3224	3231	rBV	76980	168992	4.58%	0.799%
20	35.525	3316	3321	3329	rBV	68508	147163	3.98%	0.696%
21	36.388	3411	3415	3421	rBV2	40955	93256	2.53%	0.441%
22	37.278	3503	3512	3531	rBV2	944832	2788267	75.50%	13.185%

Sum of corrected areas: 21146711

LB.D NP091101.M Wed Oct 03 10:09:37 2001

LSC Report - Integrated Chromatogram

File : C:\HPCHEM\1\DATA\OCT0201\LB.D
Operator : 209 GALOS
Acquired : 2 Oct 2001 3:06 pm using AcqMethod NP091101
Instrument : GC/MS 209
Sample Name: 9/20
Misc Info :
Vial Number: 2
Quant File :NP091101.RES (RTE Integrator)



LB.D NP091101.M Wed Oct 03 10:09:40 2001

Library Search Compound Report

Data File : C:\HPCHEM\1\DATA\OCT0201\LB.D
Acq On : 2 Oct 2001 3:06 pm
Sample : 9/20
Misc :
MS Integration Params: LSCINT.P

Vial: 2
Operator: 209 GALOS
Inst : GC/MS 209
Multiplr: 1.00

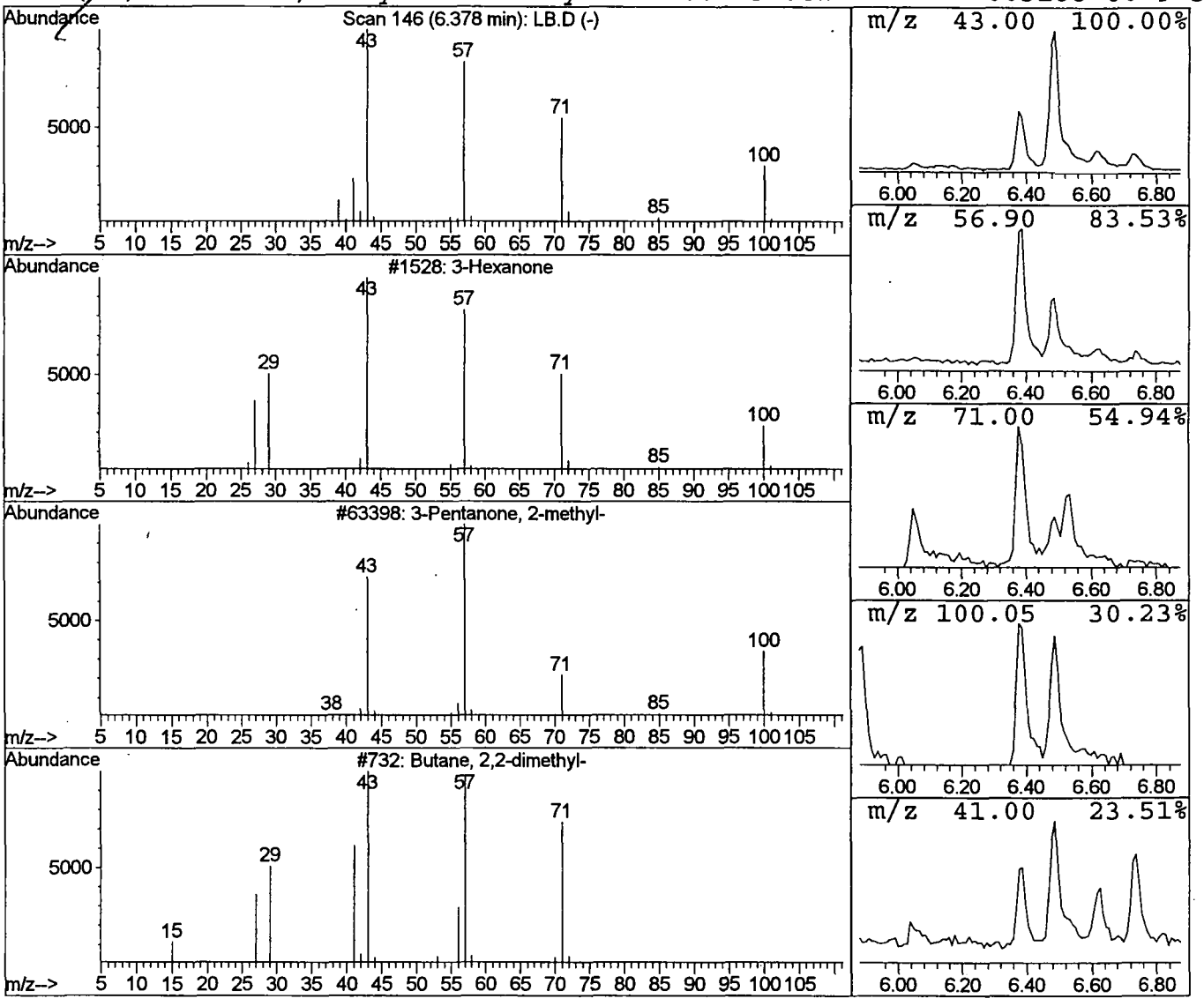
Quant Method : C:\HPCHEM\1\METHODS\NP091101.M (RTE Integrator)
Title : 209 625/TCLP calibration table 7/16/01
Library : C:\DATABASE\NBS75K.L

****

Peak Number 1 3-Hexanone Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.38	0.60 ug/l	76825	1,4-Dichlorobenzene-d4	11.78

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	3-Hexanone	100	C6H12O	000589-38-8	91
2	3-Pentanone, 2-methyl-	100	C6H12O	000565-69-5	47
3	Butane, 2,2-dimethyl-	86	C6H14	000075-83-2	36
4	3(2H)-Furanone, dihydro-2-methyl-	100	C5H8O2	003188-00-9	35



Library Search Compound Report

Data File : C:\HPCHEM\1\DATA\OCT0201\LB.D
Acq On : 2 Oct 2001 3:06 pm
Sample : 9/20
Misc :
MS Integration Params: LSCINT.P

Vial: 2
Operator: 209 GALOS
Inst : GC/MS 209
Multiplr: 1.00

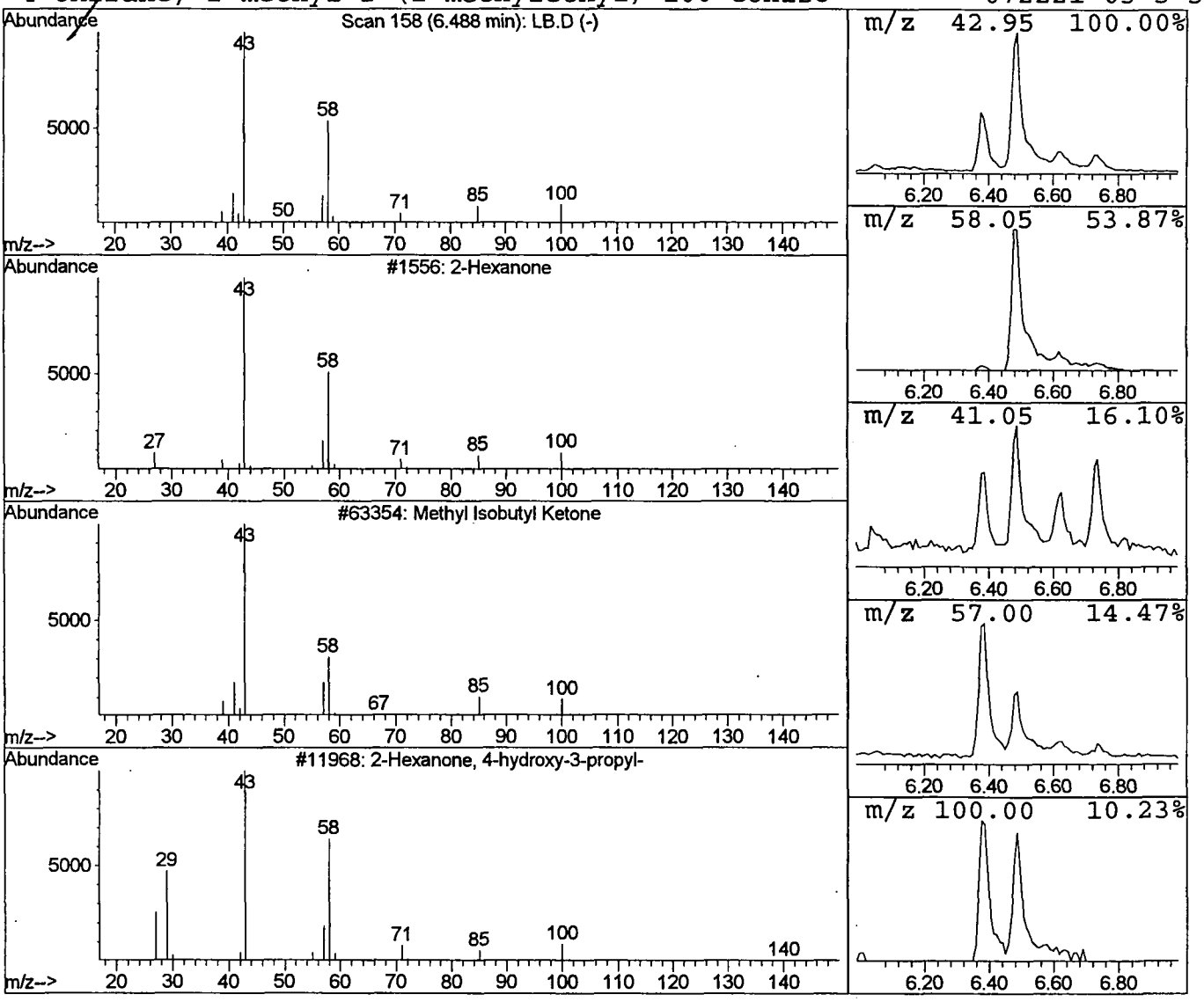
Quant Method : C:\HPCHEM\1\METHODS\NP091101.M (RTE Integrator)
Title : 209 625/TCLP calibration table 7/16/01
Library : C:\DATABASE\NBS75K.L

****

Peak Number 2 2-Hexanone Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.49	0.83 ug/l	107024	1,4-Dichlorobenzene-d4	11.78

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Hexanone	100	C6H12O	000591-78-6	91
2	Methyl Isobutyl Ketone	100	C6H12O	000108-10-1	58
3	2-Hexanone, 4-hydroxy-3-propyl-	158	C9H18O2	062338-17-4	56
4	Oxirane, 2-methyl-2-(1-methylethyl)	100	C6H12O	072221-03-5	50



Library Search Compound Report

Data File : C:\HPCHEM\1\DATA\OCT0201\LB.D
Acq On : 2 Oct 2001 3:06 pm
Sample : 9/20
Misc :
MS Integration Params: LSCINT.P

Vial: 2
Operator: 209 GALOS
Inst : GC/MS 209
Multiplr: 1.00

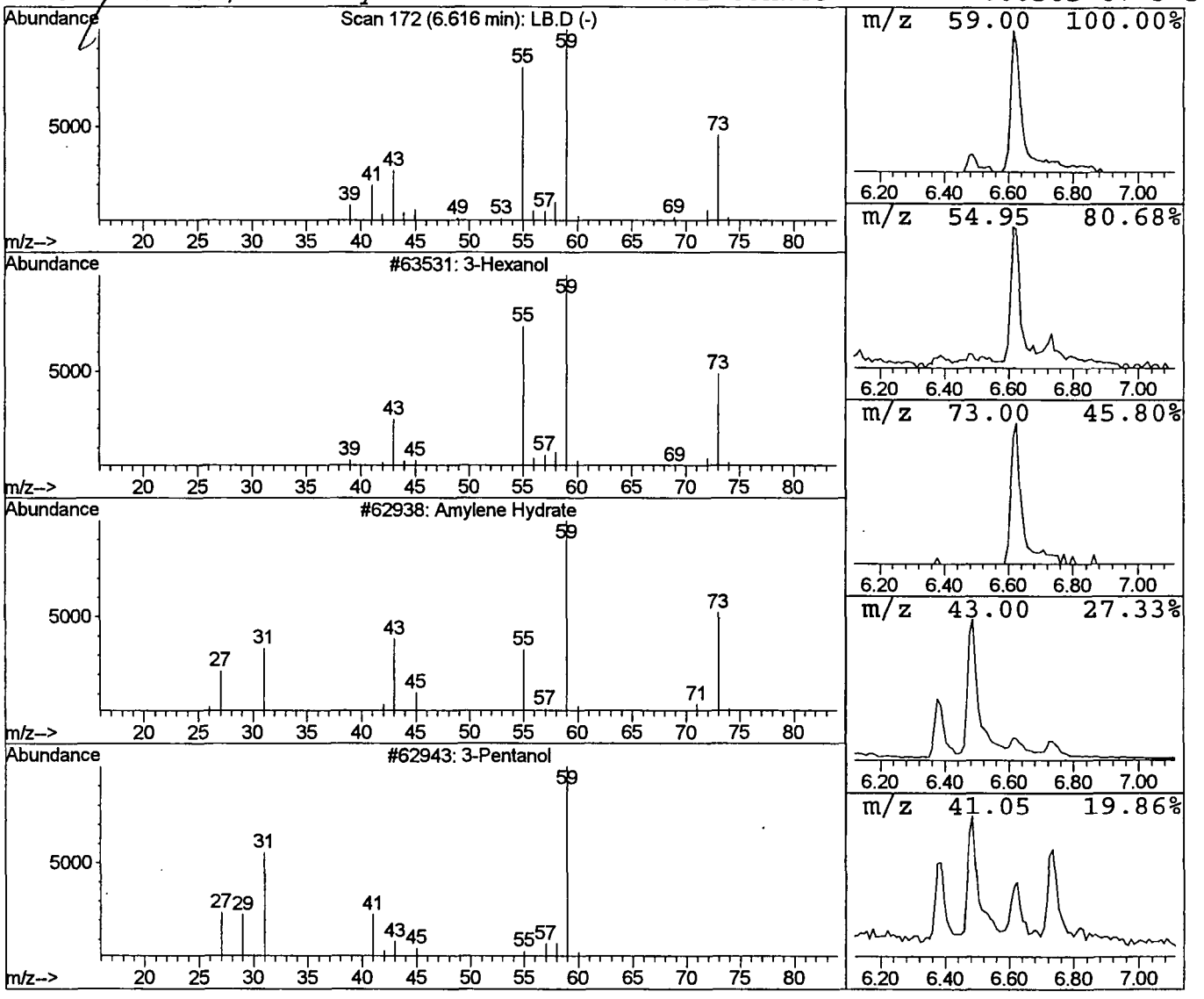
Quant Method : C:\HPCHEM\1\METHODS\NP091101.M (RTE Integrator)
Title : 209 625/TCLP calibration table 7/16/01
Library : C:\DATABASE\NBS75K.L

****

Peak Number 3 3-Hexanol Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.62	0.41 ug/l	53199	1,4-Dichlorobenzene-d4	11.78

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	3-Hexanol	102	C6H14O	000623-37-0	90
2	Amylene Hydrate	88	C5H12O	000075-85-4	45
3	3-Pentanol	88	C5H12O	000584-02-1	45
4	3-Pentanol, 2-methyl-	102	C6H14O	000565-67-3	36



Library Search Compound Report

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Acq On : 2 Oct 2001 3:06 pm
Sample : 9/20
Misc :
MS Integration Params: LSCINT.P

Vial: 2
Operator: 209 GALOS
Inst : GC/MS 209
Multiplr: 1.00

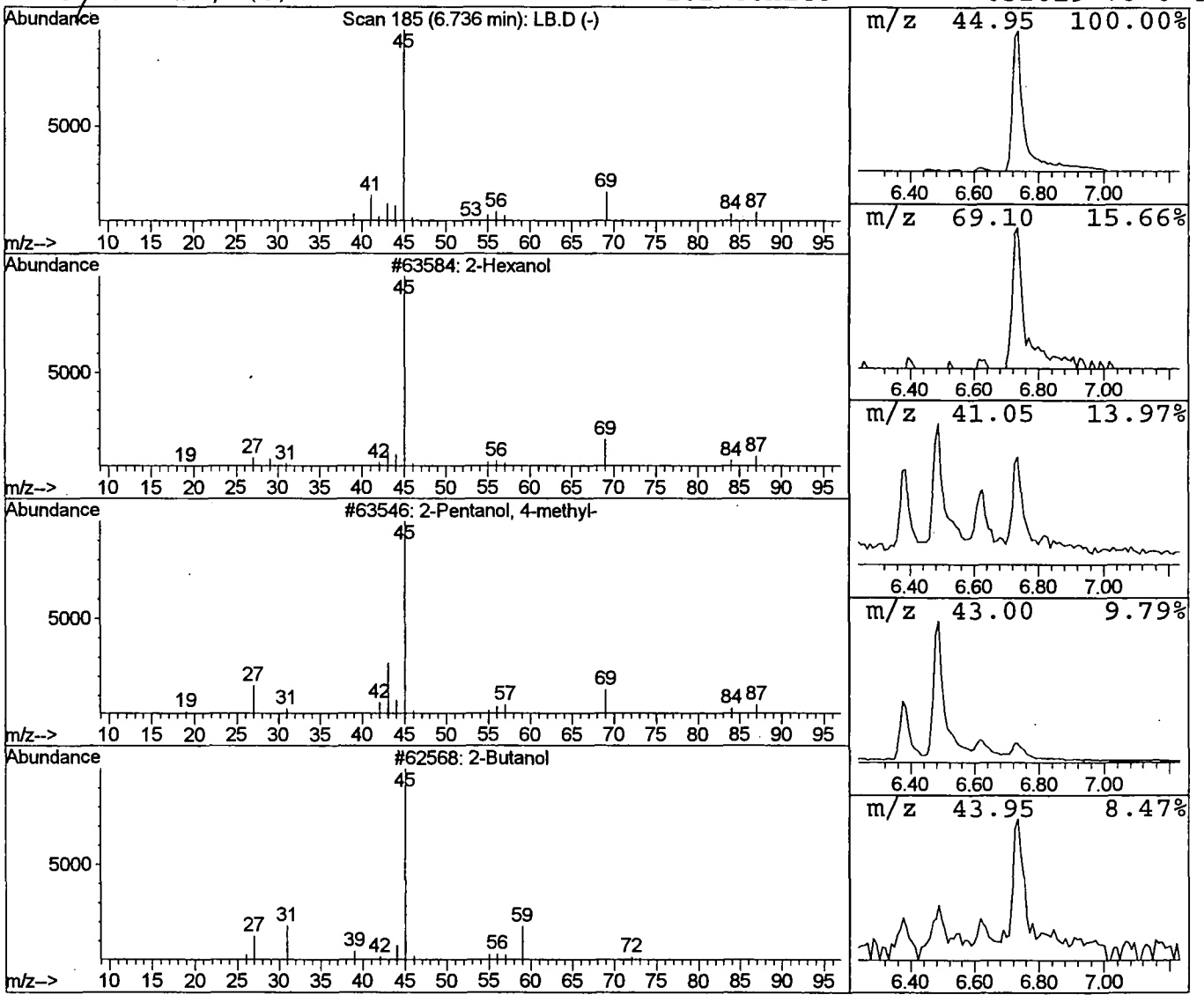
Quant Method : C:\HPCHEM\1\METHODS\NP091101.M (RTE Integrator)
Title : 209 625/TCLP calibration table 7/16/01
Library : C:\DATABASE\NBS75K.L

****

Peak Number 4 2-Hexanol Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.74	0.69 ug/l	88381	1,4-Dichlorobenzene-d4	11.78

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Hexanol	102	C6H14O	000626-93-7	83
2	2-Pentanol, 4-methyl-	102	C6H14O	000108-11-2	78
3	2-Butanol	74	C4H10O	000078-92-2	50
4	2-Hexanol, (S)-	102	C6H14O	052019-78-0	43



Library Search Compound Report

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Acq On : 2 Oct 2001 3:06 pm
Sample : 9/20
Misc :
MS Integration Params: LSCINT.P

Vial: 2
Operator: 209 GALOS
Inst : GC/MS 209
Multiplr: 1.00

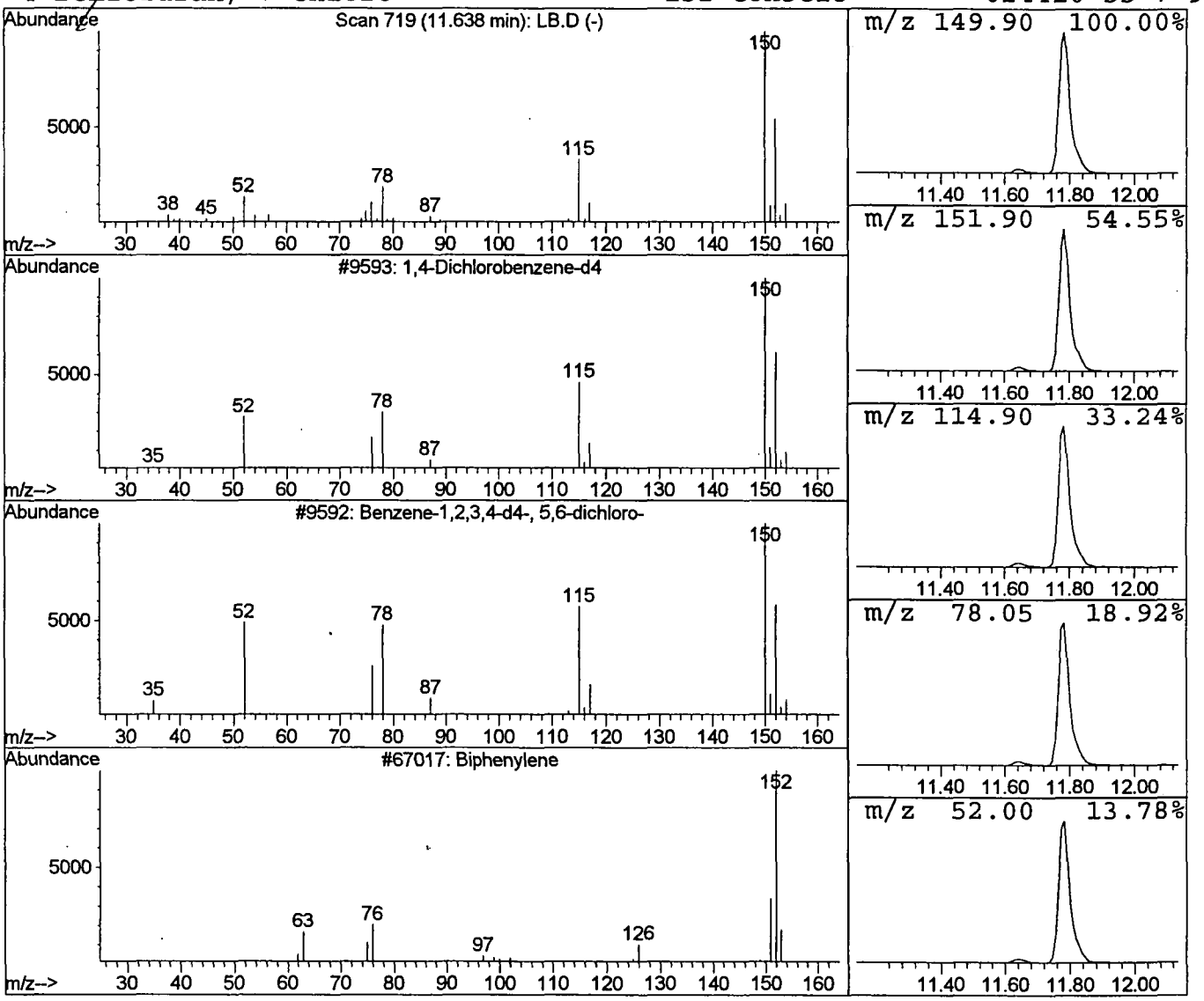
Quant Method : C:\HPCHEM\1\METHODS\NP091101.M (RTE Integrator)
Title : 209 625/TCLP calibration table 7/16/01
Library : C:\DATABASE\NBS75K.L

****

Peak Number 5 1,4-Dichlorobenzene-d4 Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.64	0.54 ug/l	69043	1,4-Dichlorobenzene-d4	11.78

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,4-Dichlorobenzene-d4	150	C6Cl2D4	003855-82-1	50
2	Benzene-1,2,3,4-d4-, 5,6-dichloro-	150	C6Cl2D4	002199-69-1	27
3	Biphenylene	152	C12H8	000259-79-0	9
4	Benzofuran, 7-chloro-	152	C8H5ClO	024410-55-7	9



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Sample : 9/20
Misc :
MS Integration Params: LSCINT.P

Vial: 2
Operator: 209 GALOS
Inst : GC/MS 209
Multiplr: 1.00

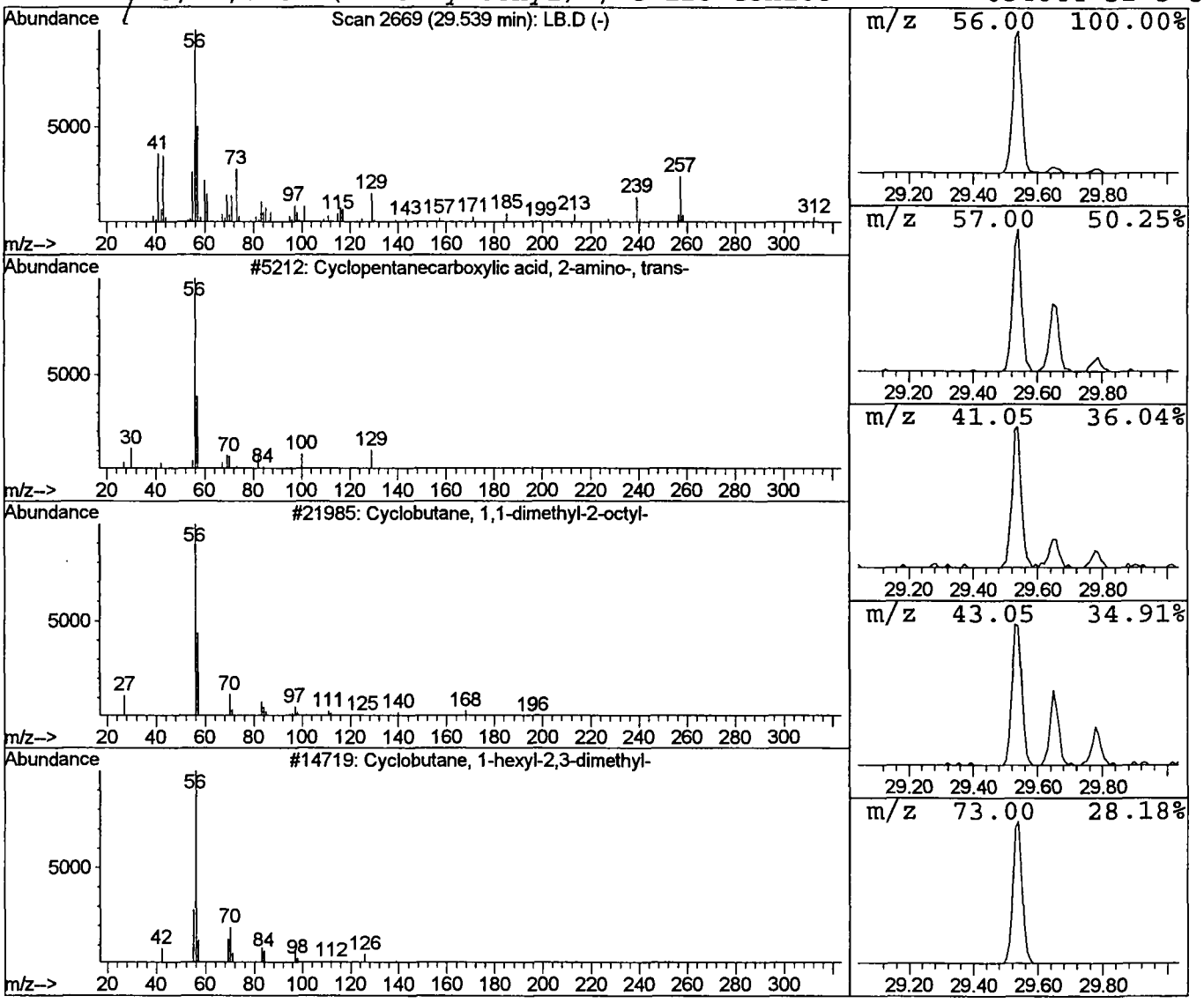
Quant Method : C:\HPCHEM\1\METHODS\NP091101.M (RTE Integrator)
Title : 209 625/TCLP calibration table 7/16/01
Library : C:\DATABASE\NBS75K.L

****

Peak Number 6 Cyclopentanecarboxylic acid, 2 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
29.54	1.71 ug/l	260921	Chrysene-d12	33.16

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclopentanecarboxylic acid, 2-amin	129	C6H11NO2	040482-05-1	38
2	Cyclobutane, 1,1-dimethyl-2-octyl-	196	C14H28	062338-30-1	35
3	Cyclobutane, 1-hexyl-2,3-dimethyl-	168	C12H24	055170-84-8	35
4	Oxirane, 2,3-bis(1-methylethyl)-, t	128	C8H16O	054644-32-5	32



Library Search Compound Report

Data File : C:\HPCHEM\1\DATA\OCT0201\LB.D
Acq On : 2 Oct 2001 3:06 pm
Sample : 9/20
Misc :
MS Integration Params: LSCINT.P

Vial: 2
Operator: 209 GALOS
Inst : GC/MS 209
Multiplr: 1.00

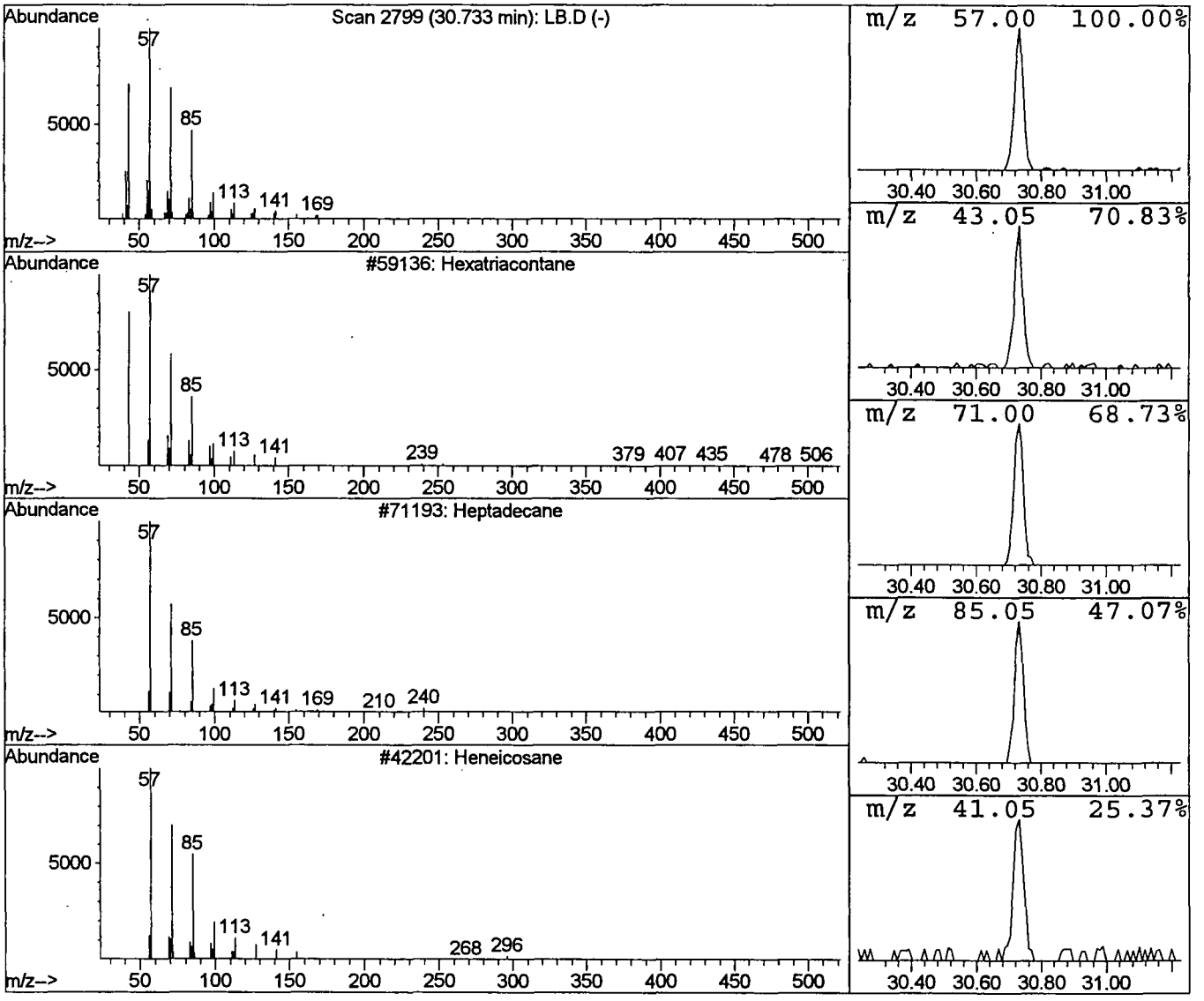
Quant Method : C:\HPCHEM\1\METHODS\NP091101.M (RTE Integrator)
Title : 209 625/TCLP calibration table 7/16/01
Library : C:\DATABASE\NBS75K.L

****

Peak Number 7 Hexatriacontane Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
30.73	0.55 ug/l	84270	Chrysene-d12	33.16

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Hexatriacontane	507	C36H74	000630-06-8	91
2	Heptadecane	240	C17H36	000629-78-7	91
3	Heneicosane	296	C21H44	000629-94-7	90
4	Tetracosane	338	C24H50	000646-31-1	90



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Sample : 9/20
Misc :
MS Integration Params: LSCINT.P

Vial: 2
Operator: 209 GALOS
Inst : GC/MS 209
Multiplr: 1.00

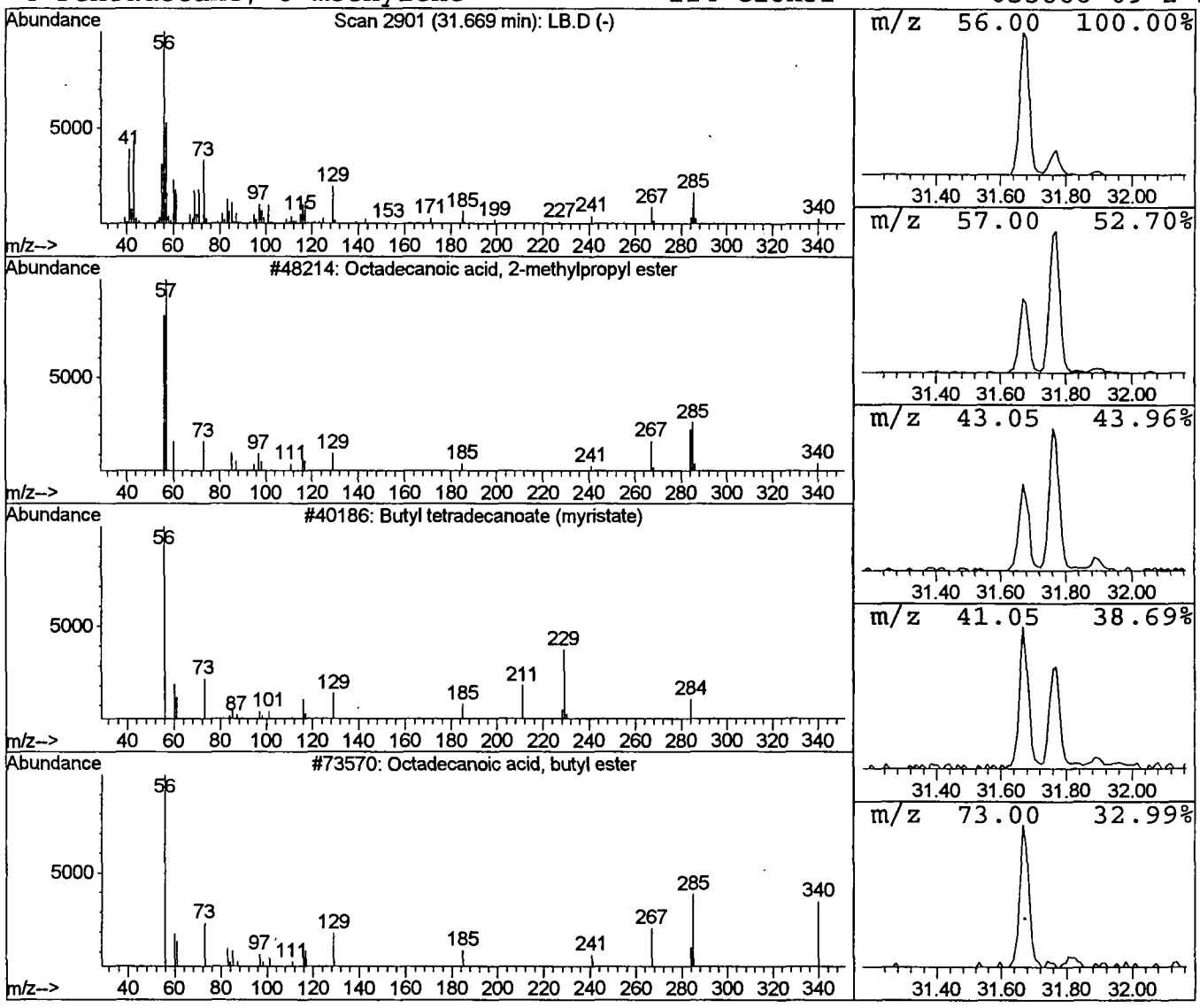
Quant Method : C:\HPCHEM\1\METHODS\NP091101.M (RTE Integrator)
Title : 209 625/TCLP calibration table 7/16/01
Library : C:\DATABASE\NBS75K.L

****

Peak Number 8 Octadecanoic acid, 2-methylpro Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
31.67	1.34 ug/l	204617	Chrysene-d12	33.16

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Octadecanoic acid, 2-methylpropyl e	340	C22H44O2	000646-13-9	87
2	Butyl tetradecanoate (myristate)	284	C18H36O2	000000-00-0	86
3	Octadecanoic acid, butyl ester	340	C22H44O2	000123-95-5	56
4	Pentadecane, 8-methylene-	224	C16H32	055668-09-2	32



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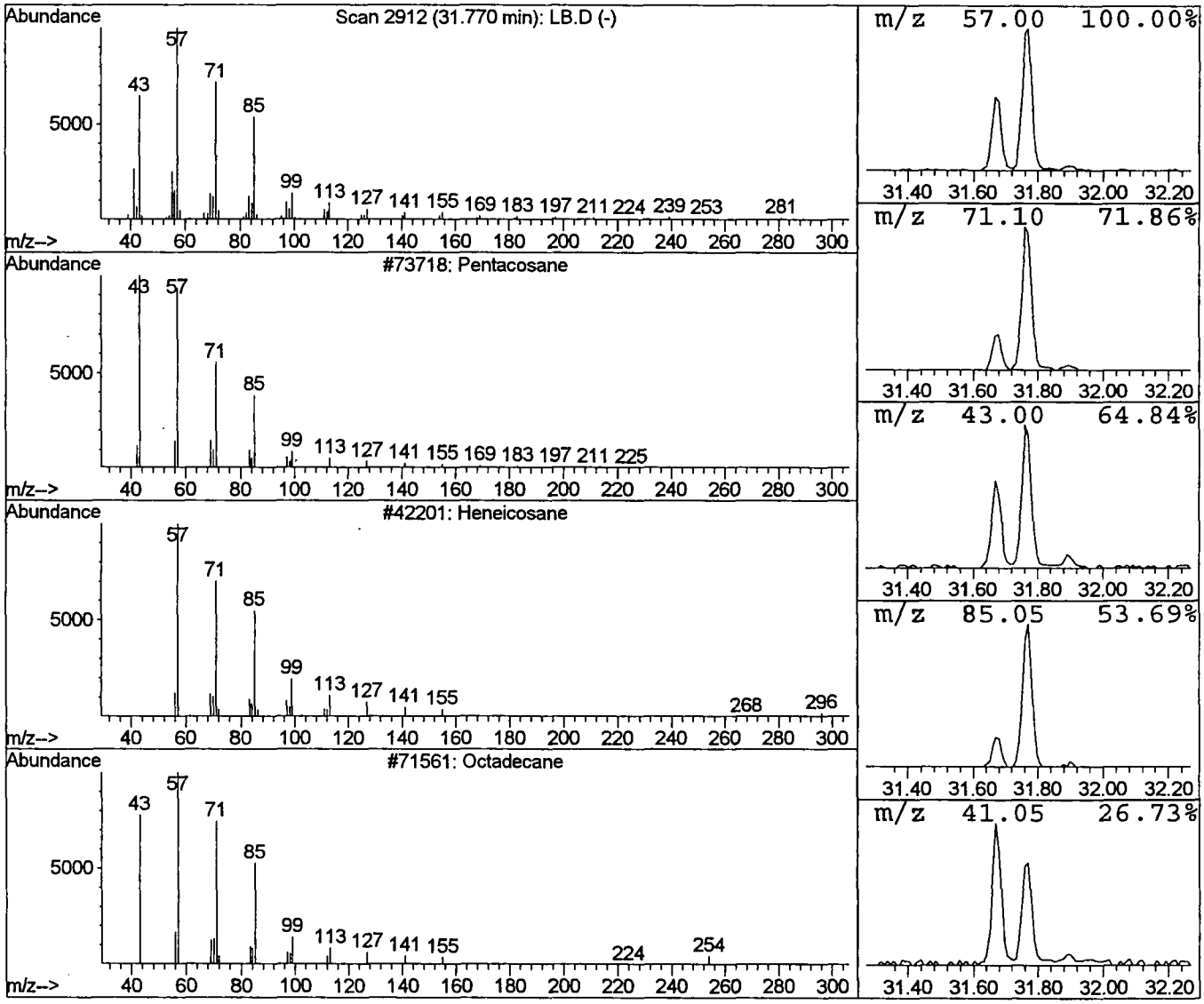
Title : 209 625/TCLP calibration table 7/16/01
Library : C:\DATABASE\NBS75K.L

****

Peak Number 9 Pentacosane Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
31.77	1.18 ug/l	179169	Chrysene-d12	33.16

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Pentacosane	352	C25H52	000629-99-2	91
2	Heneicosane	296	C21H44	000629-94-7	91
3	Octadecane	254	C18H38	000593-45-3	91
4	Tetracosane	338	C24H50	000646-31-1	90



Library Search Compound Report

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Sample : 9/20
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MS Integration Params: LSCINT.P

Vial: 2
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Inst : GC/MS 209
Multiplr: 1.00

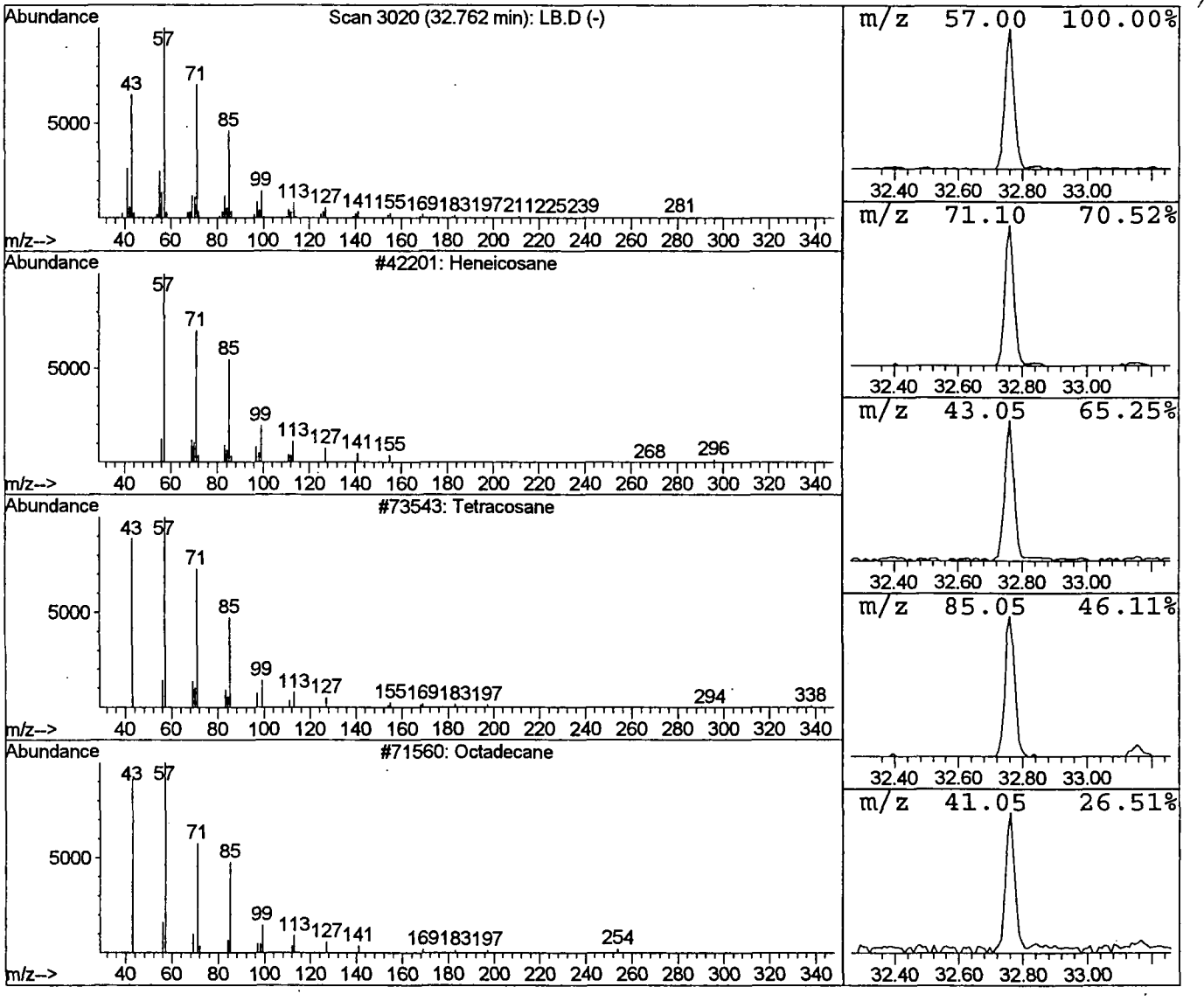
Quant Method : C:\HPCHEM\1\METHODS\NP091101.M (RTE Integrator)
Title : 209 625/TCLP calibration table 7/16/01
Library : C:\DATABASE\NBS75K.L

****

Peak Number 10 Heneicosane Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
32.76	1.20 ug/l	183463	Chrysene-d12	33.16

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heneicosane	296	C21H44	000629-94-7	90
2	Tetracosane	338	C24H50	000646-31-1	90
3	Octadecane	254	C18H38	000593-45-3	90
4	Tricosane	324	C23H48	000638-67-5	87



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Quant Method : C:\HPCHEM\1\METHODS\NP091101.M (RTE Integrator)
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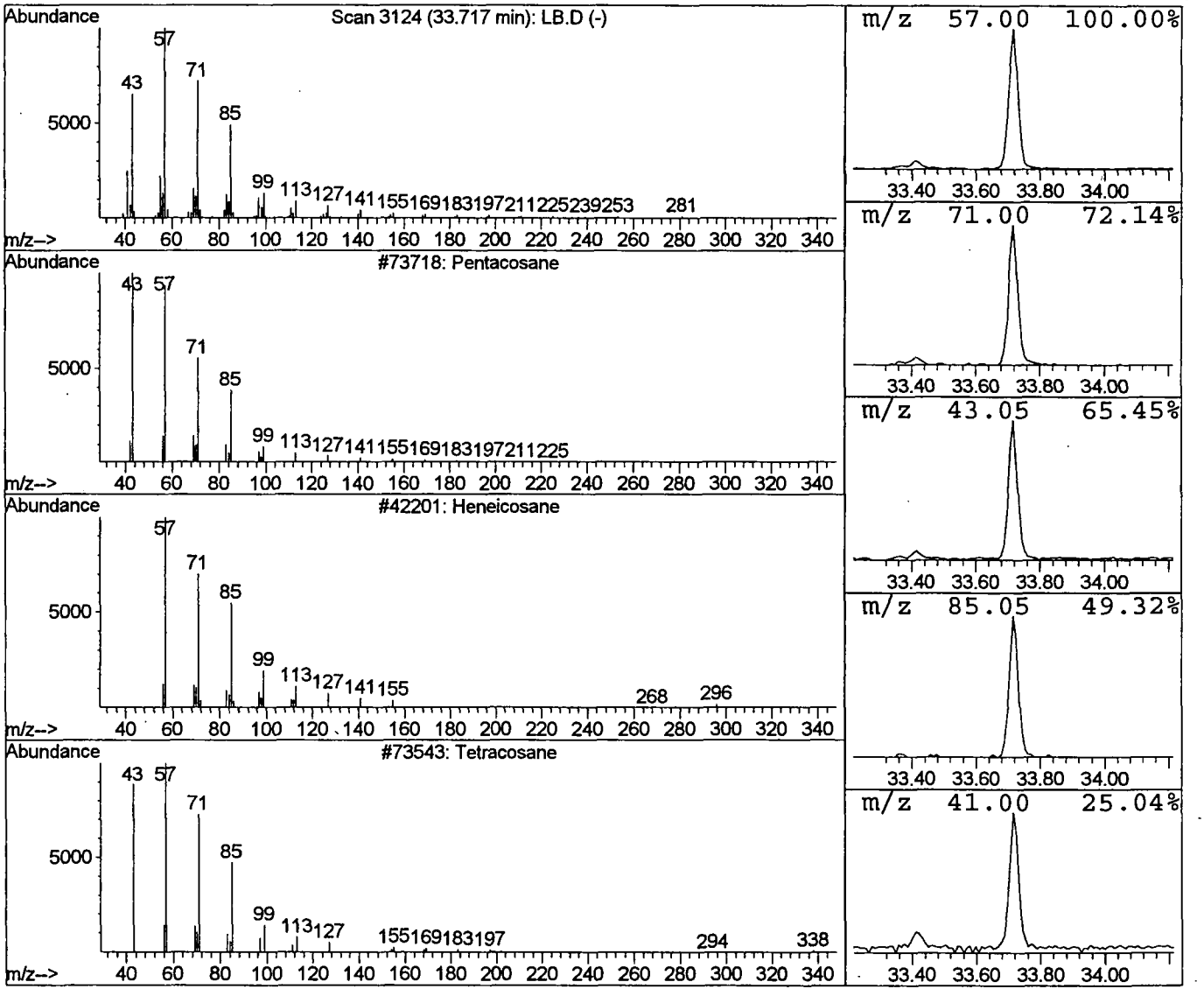
****

Peak Number 11 Pentacosane — Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
33.72	1.45 ug/l	221588	Chrysene-d12	33.16

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Pentacosane	352	C25H52	000629-99-2	91
2	Heneicosane	296	C21H44	000629-94-7	91
3	Tetracosane	338	C24H50	000646-31-1	91
4	Tricosane	324	C23H48	000638-67-5	91



Library Search Compound Report

Data File : C:\HPCHEM\1\DATA\OCT0201\LB.D
Acq On : 2 Oct 2001 3:06 pm
Sample : 9/20
Misc :
MS Integration Params: LSCINT.P

Vial: 2
Operator: 209 GALOS
Inst : GC/MS 209
Multiplr: 1.00

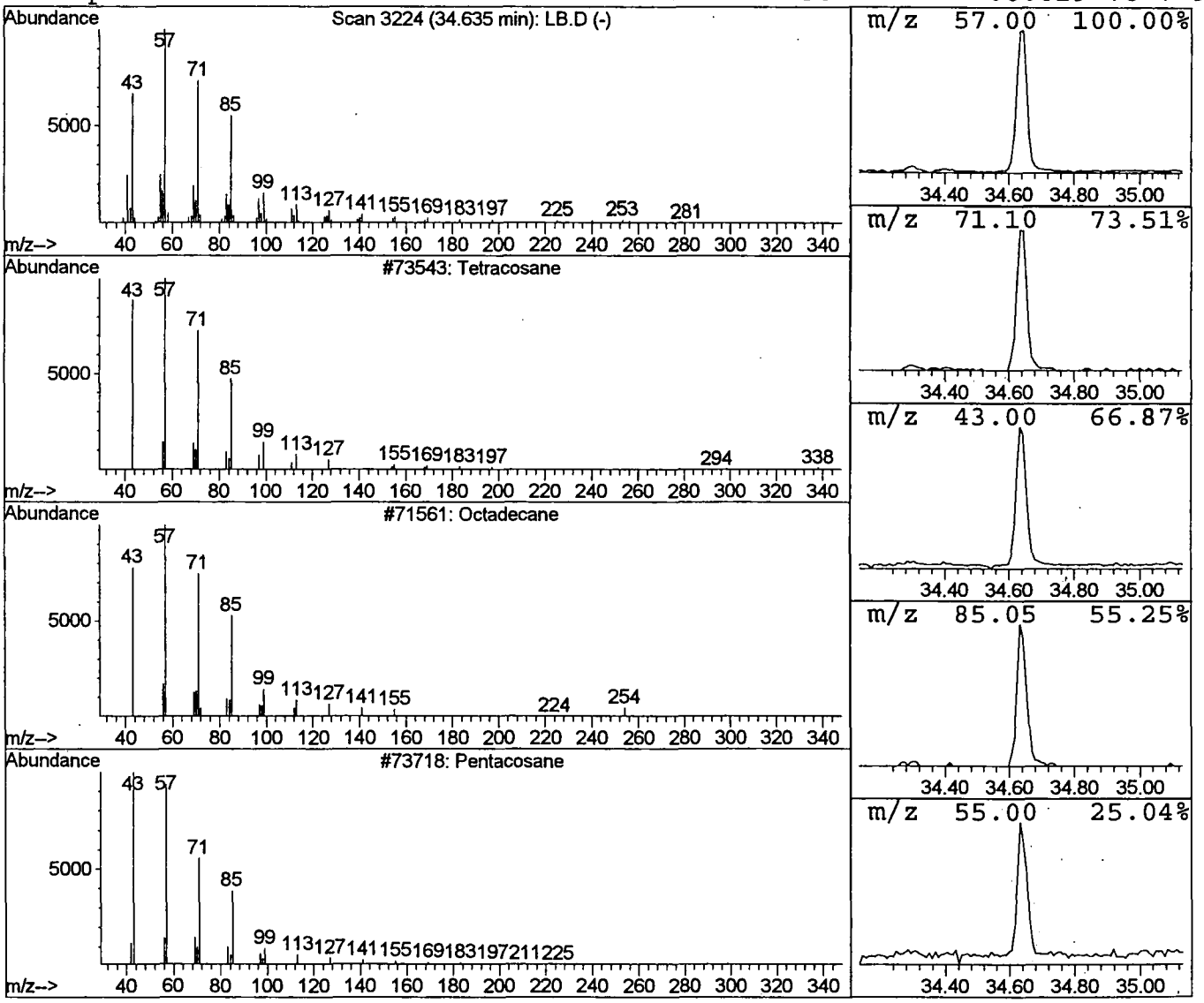
Quant Method : C:\HPCHEM\1\METHODS\NP091101.M (RTE Integrator)
Title : 209 625/TCLP calibration table 7/16/01
Library : C:\DATABASE\NBS75K.L

****

Peak Number 12 Tetracosane Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
34.63	1.11 ug/l	168992	Chrysene-d12	33.16

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Tetracosane	338	C24H50	000646-31-1	91
2	Octadecane	254	C18H38	000593-45-3	91
3	Pentacosane	352	C25H52	000629-99-2	91
4	Heptadecane	240	C17H36	000629-78-7	91



Library Search Compound Report

Data File : C:\HPCHEM\1\DATA\OCT0201\LB.D
Acq On : 2 Oct 2001 3:06 pm
Sample : 9/20
Misc :
MS Integration Params: LSCINT.P

Vial: 2
Operator: 209 GALOS
Inst : GC/MS 209
Multiplr: 1.00

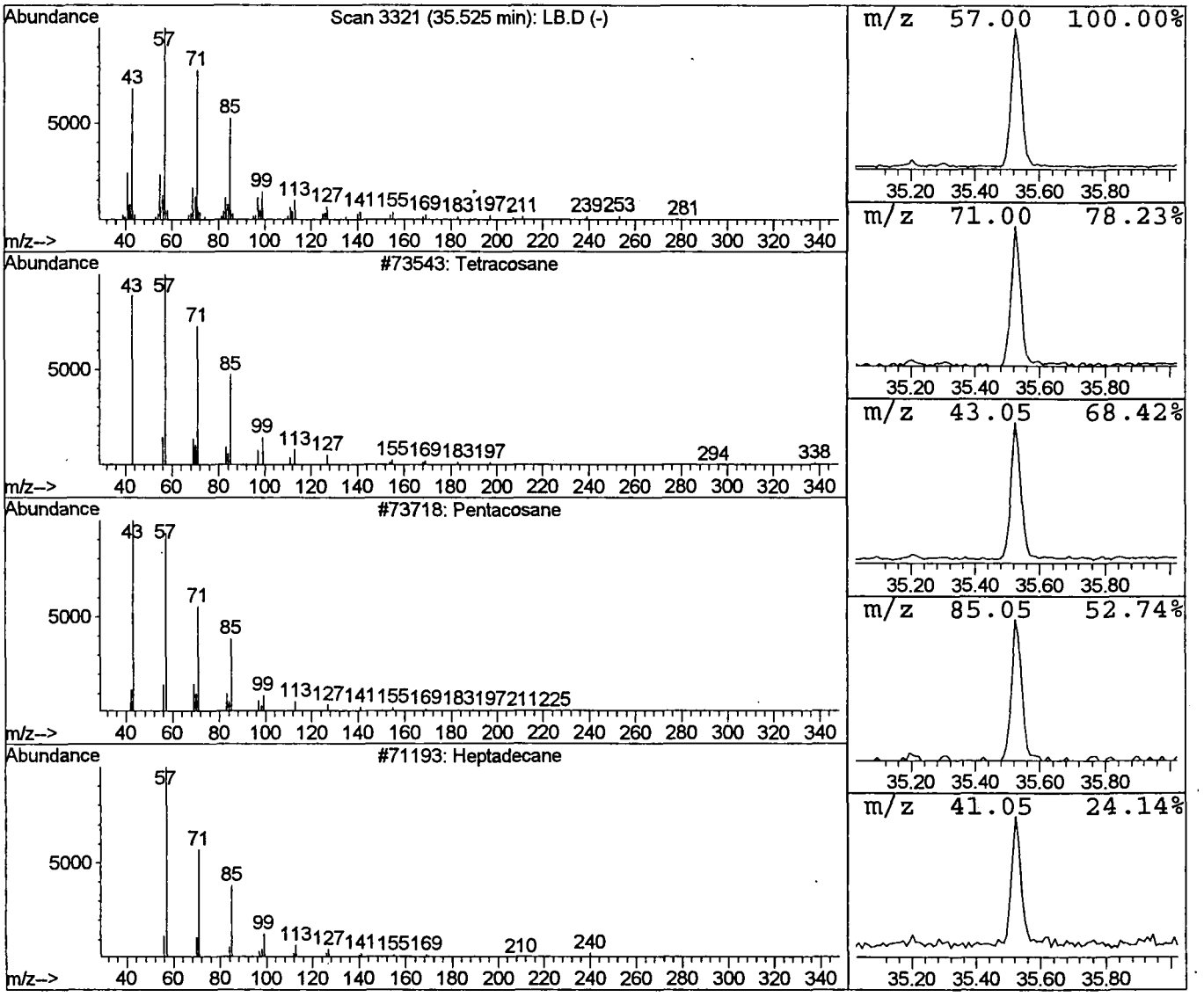
Quant Method : C:\HPCHEM\1\METHODS\NP091101.M (RTE Integrator)
Title : 209 625/TCLP calibration table 7/16/01
Library : C:\DATABASE\NBS75K.L

****

Peak Number 13 Tetracosane Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
35.53	1.06 ug/l	147163	Perylene-d12	37.28

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Tetracosane	338	C24H50	000646-31-1	91
2	Pentacosane	352	C25H52	000629-99-2	91
3	Heptadecane	240	C17H36	000629-78-7	91
4	Hexatriacontane	507	C36H74	000630-06-8	91



Library Search Compound Report

Data File : C:\HPCHEM\1\DATA\OCT0201\LB.D
Acq On : 2 Oct 2001 3:06 pm
Sample : 9/20
Misc :
MS Integration Params: LSCINT.P

Vial: 2
Operator: 209 GALOS
Inst : GC/MS 209
Multiplr: 1.00

Quant Method : C:\HPCHEM\1\METHODS\NP091101.M (RTE Integrator)
Title : 209 625/TCLP calibration table 7/16/01
Library : C:\DATABASE\NBS75K.L

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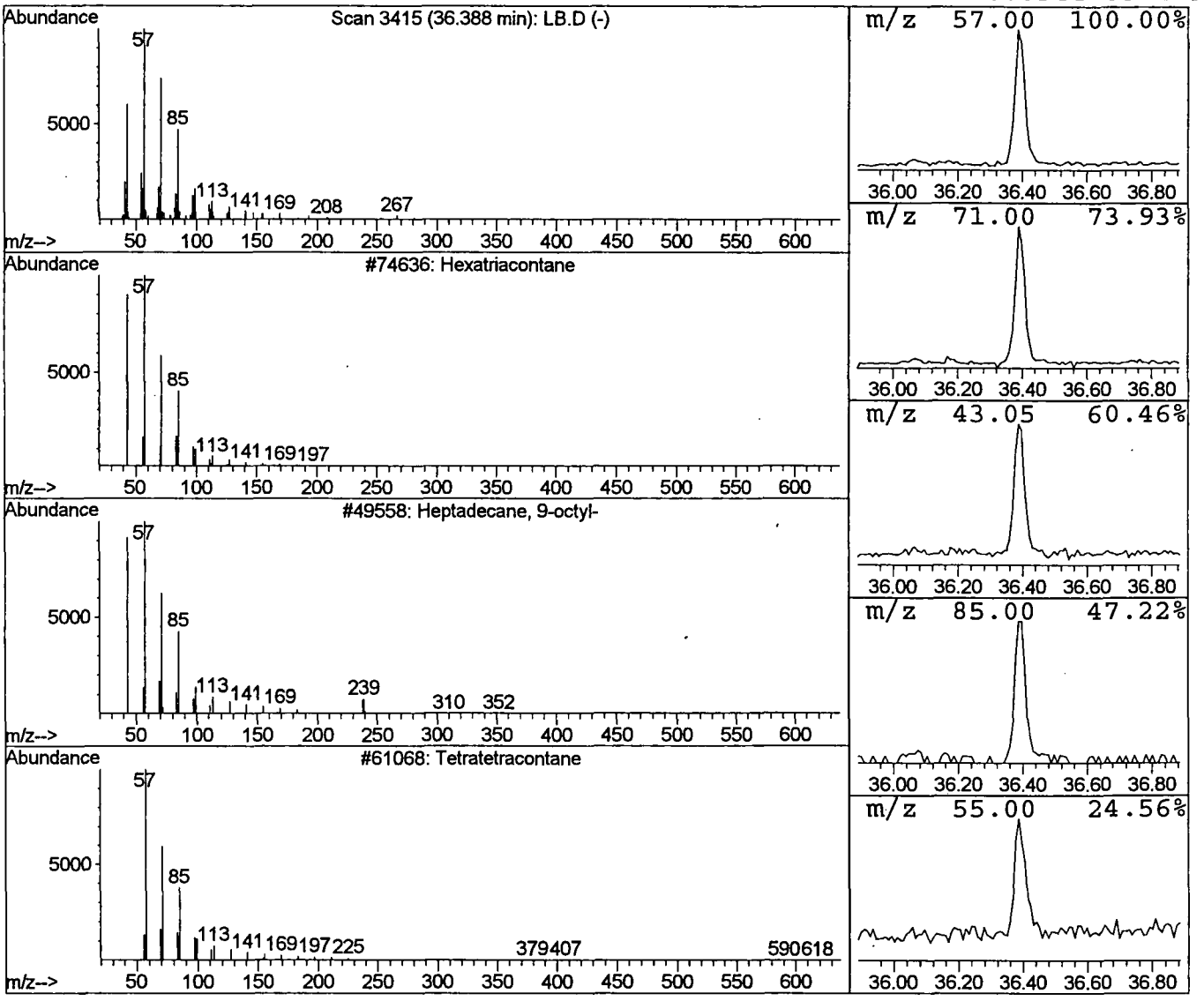
Peak Number 14 Hexatriacontane Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
36.39	0.67 ug/l	93256	Perylene-d12	37.28

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Hexatriacontane	507	C36H74	000630-06-8	91
2	Heptadecane, 9-octyl-	352	C25H52	007225-64-1	91
3	Tetratetracontane	619	C44H90	007098-22-8	91
4	Dotriacontane	451	C32H66	000544-85-4	91



Tentatively Identified Compound (LSC) summary

Operator ID: 209 GALOS Date Acquired: 2 Oct 2001 3:06 pm
Data File: C:\HPCHEM\1\DATA\OCT0201\LB.D
Name: 9/20
Misc:
Method: C:\HPCHEM\1\METHODS\NP091101.M (RTE Integrator)
Title: 209 625/TCLP calibration table 7/16/01
Library Searched: C:\DATABASE\NBS75K.L

TIC Top Hit name	RT	EstConc	Units	Area	IntStd	ISRT	ISArea	ISConc
3-Hexanone	6.38	0.6	ug/l	76825	ISTD01	11.78	2566400	20.0
2-Hexanone	6.49	0.8	ug/l	107024	ISTD01	11.78	2566400	20.0
3-Hexanol	6.62	0.4	ug/l	53199	ISTD01	11.78	2566400	20.0
2-Hexanol	6.74	0.7	ug/l	88381	ISTD01	11.78	2566400	20.0
1,4-Dichlorobenzene-	11.64	0.5	ug/l	69043	ISTD01	11.78	2566400	20.0
Cyclopentanecarboxyl	29.54	1.7	ug/l	260921	ISTD05	33.16	3048400	20.0
Hexatriacontane	30.73	0.6	ug/l	84270	ISTD05	33.16	3048400	20.0
Octadecanoic acid, 2	31.67	1.3	ug/l	204617	ISTD05	33.16	3048400	20.0
Pentacosane	31.77	1.2	ug/l	179169	ISTD05	33.16	3048400	20.0
Heneicosane	32.76	1.2	ug/l	183463	ISTD05	33.16	3048400	20.0
Pentacosane	33.72	1.5	ug/l	221588	ISTD05	33.16	3048400	20.0
Tetracosane	34.63	1.1	ug/l	168992	ISTD05	33.16	3048400	20.0
Tetracosane	35.53	1.1	ug/l	147163	ISTD06	37.28	2788270	20.0
Hexatriacontane	36.39	0.7	ug/l	93256	ISTD06	37.28	2788270	20.0

LB.D NP091101.M Wed Oct 03 10:10:14 2001