User: Galos, Marie
Westchester Dept. of Labs & Research
Date: 10/16/2001 Time: 09:17:18

Sample: AD23471
Analysis: $GCMS GCMS Scan For Unknowns
Units: ug
Started: 10/16/01 09:15 Ended: 10/16/01 09:15 Analyst: MG
Page: 1

	Component Name	Vial	Result
1	Other unknown compounds		see comment

Quantitation Report (QT Reviewed)

Data File : C:\HPCHEM\1\DATA\OCT0901\23471.D Vial: 19
Acq On : 10 Oct 2001 6:53 am Operator: 209 GALOS
Sample : gc/ms unknown Inst : GC/MS 209
Misc : Multiplr: 1.00
MS Integration Params: RTEINT.P
Quant Time: Oct 11 12:23 2001 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

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	11.76	152	363156	20.00	ug/l	-0.14
19) Naphthalene-d8	15.37	136	1432130	20.00	ug/l	-0.15
31) Acenaphthene-d10	20.66	164	660057	20.00	ug/l	-0.15
49) Phenanthrene-d10	25.08	188	901887	20.00	ug/l	-0.15
59) Chrysene-d12	33.12	240	548238	20.00	ug/l	-0.16
68) Perylene-d12	37.23	264	375117	20.00	ug/l	-0.19
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	11.75	132	181	0.01	ug/l	0.34
Spiked Amount 75.000			Recovery =	0.01%		
7) 1,2-Dichlorobenzene-d4	12.28	152	3850	0.22	ug/l	-0.15
Spiked Amount 50.000			Recovery =	0.44%		
20) Nitrobenzene-d5	0.00	82	0	0.00	ug/l	
Spiked Amount 50.000			Recovery =	0.00%		
32) 2-Fluorobiphenyl	18.80	172	1447	0.03	ug/l	-0.02
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	0.00	79	0	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 (QT Reviewed)

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Last Update : Thu Sep 13 12:47:04 2001
Response via : Initial Calibration
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	Compound	R.T.	QIon	Response	Conc Unit	Qvalue
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-Chlorophenyl-phenylether	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	168	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	27.08	149	2646	N.D.	
58)	Fluoranthene	0.00	202	0	N.D.	
60)	3,3'-Dichlorobenzidine	0.00	252	0	N.D.	
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	33.12	228	1388	N.D.	
66)	Bis(2-ethylhexyl)phthalate	33.38	149	2035	N.D.	
67)	Chrysene	33.12	228	1388	N.D.	
69)	Di-n-Octylphthalate	0.00	149	0	N.D.	
70)	Benzo(b)fluoranthene	0.00	252	0	N.D.	

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MS Integration Params: RTEINT.P
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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

Compound	R.T.	QIon	Response	Conc Unit	Qvalue
71) Benzo(k)fluoranthene	0.00	252	0	N.D.	
72) Benzo(a)pyrene	37.23	252	1534	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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Quantitation Report

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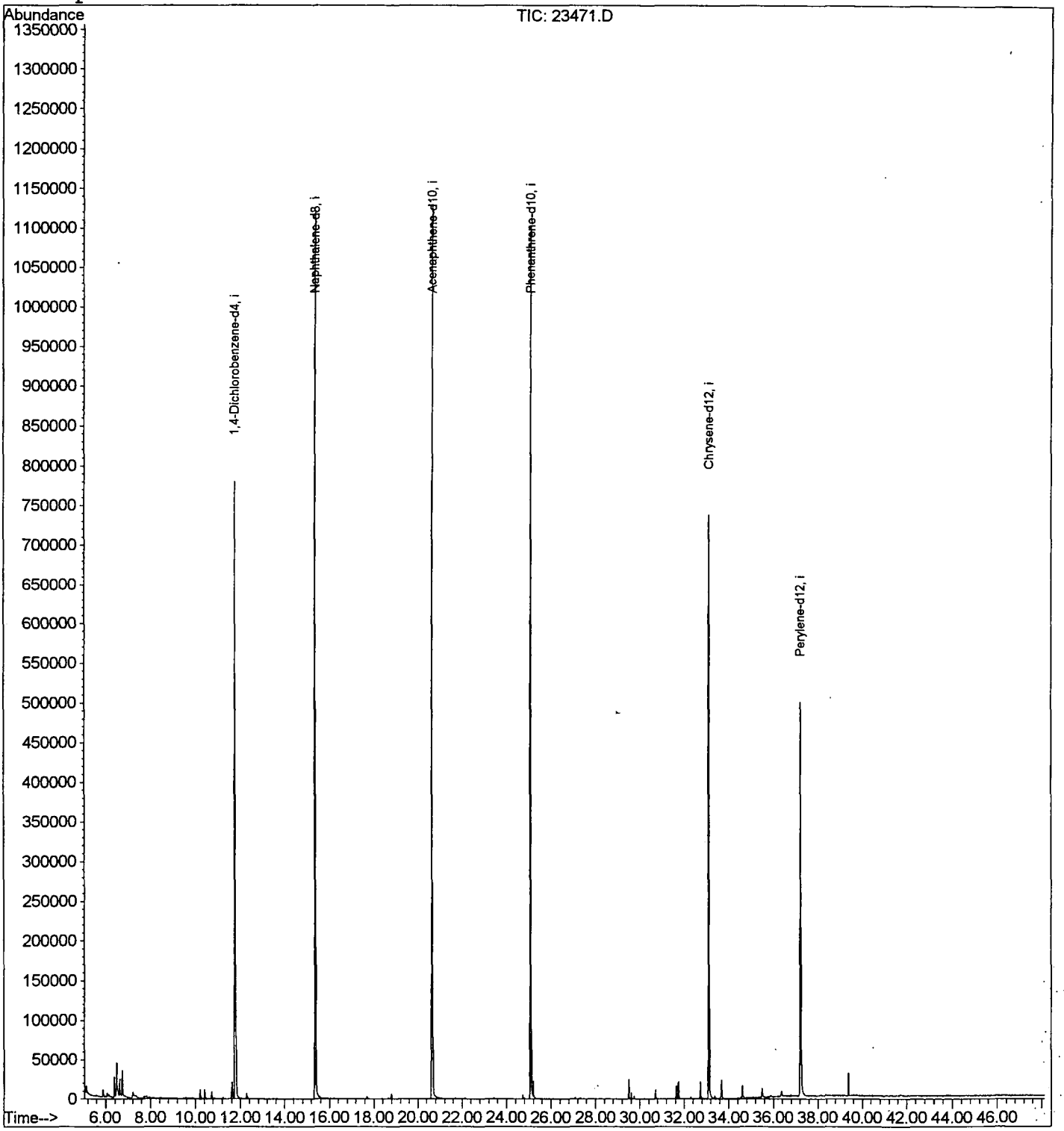
Title : 209 625/TCLP calibration table 7/16/01
Last Update : Thu Sep 13 12:47:04 2001
Response via : Initial Calibration



LSC Area Percent Report

Data File : C:\HPCHEM\1\DATA\OCT0901\23471.D Vial: 19
Acq On : 10 Oct 2001 6:53 am Operator: 209 GALOS
Sample : gc/ms unknown Inst : GC/MS 209
Misc : Multiplr: 1.00
MS Integration Params: LSCINT.P

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

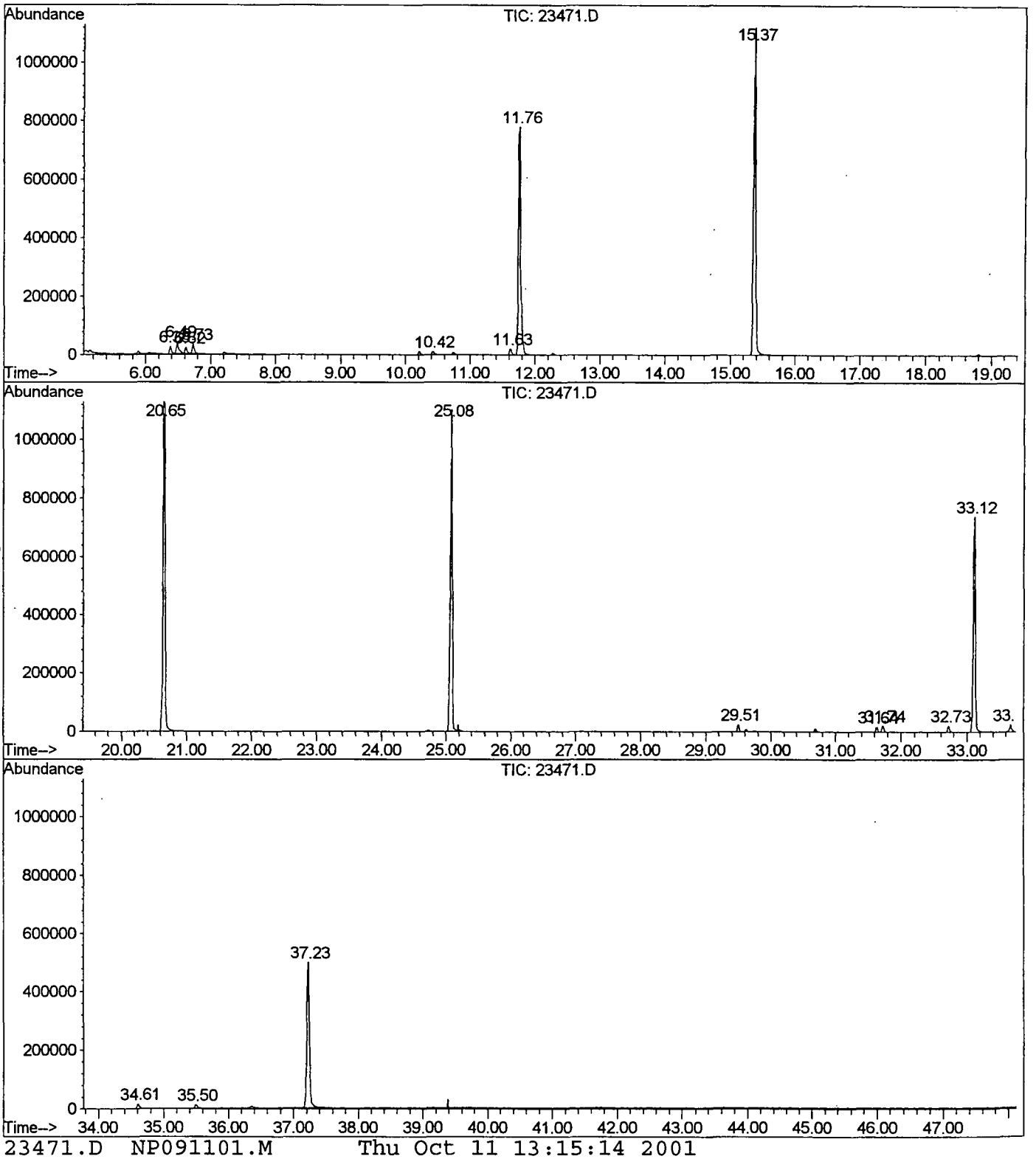
peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	peak area	peak % max.	% of total
1	6.384	142	146	151	rBV	25926	51102	1.93%	0.385%
2	6.485	153	157	168	rVV	43057	110113	4.17%	0.830%
3	6.623	168	172	180	rVV	22015	48469	1.83%	0.365%
4	6.733	180	184	193	rVB	32458	65171	2.47%	0.491%
5	10.424	582	586	595	rBV	11672	29084	1.10%	0.219%
6	11.626	712	717	726	rVB	21042	50861	1.92%	0.383%
7	11.764	726	732	749	rBV	780100	1913976	72.41%	14.425%
8	15.372	1118	1125	1142	rBV	1121453	2622393	99.21%	19.764%
9	20.650	1693	1700	1714	rBV	1129604	2643166	100.00%	19.921%
10	25.085	2175	2183	2194	rBV	1102122	2342156	88.61%	17.652%
11	29.509	2660	2665	2670	rVB	24461	44346	1.68%	0.334%
12	31.639	2889	2897	2903	rBV	16601	35046	1.33%	0.264%
13	31.740	2903	2908	2913	rVB	21527	44576	1.69%	0.336%
14	32.732	3010	3016	3021	rBV	21464	44327	1.68%	0.334%
15	33.117	3051	3058	3074	rBV2	737905	1713808	64.84%	12.916%
16	33.687	3108	3120	3127	rBV2	23747	55225	2.09%	0.416%
17	34.614	3216	3221	3226	rBV	15356	32192	1.22%	0.243%
18	35.495	3311	3317	3322	rBV2	11603	28532	1.08%	0.215%
19	37.230	3498	3506	3523	rBV2	497255	1393947	52.74%	10.506%

Sum of corrected areas: 13268490

23471.D NP091101.M Thu Oct 11 13:15:12 2001

LSC Report - Integrated Chromatogram

File : C:\HPCHEM\1\DATA\OCT0901\23471.D
Operator : 209 GALOS
Acquired : 10 Oct 2001 6:53 am using AcqMethod NP091101
Instrument : GC/MS 209
Sample Name: gc/ms unknown
Misc Info :
Vial Number: 19
Quant File :NP091101.RES (RTE Integrator)



23471.D NP091101.M Thu Oct 11 13:15:14 2001

Library Search Compound Report

Data File : C:\HPCHEM\1\DATA\OCT0901\23471.D
Acq On : 10 Oct 2001 6:53 am
Sample : gc/ms unknown
Misc :
MS Integration Params: LSCINT.P

Vial: 19
Operator: 209 GALOS
Inst : GC/MS 209
Multiplr: 1.00

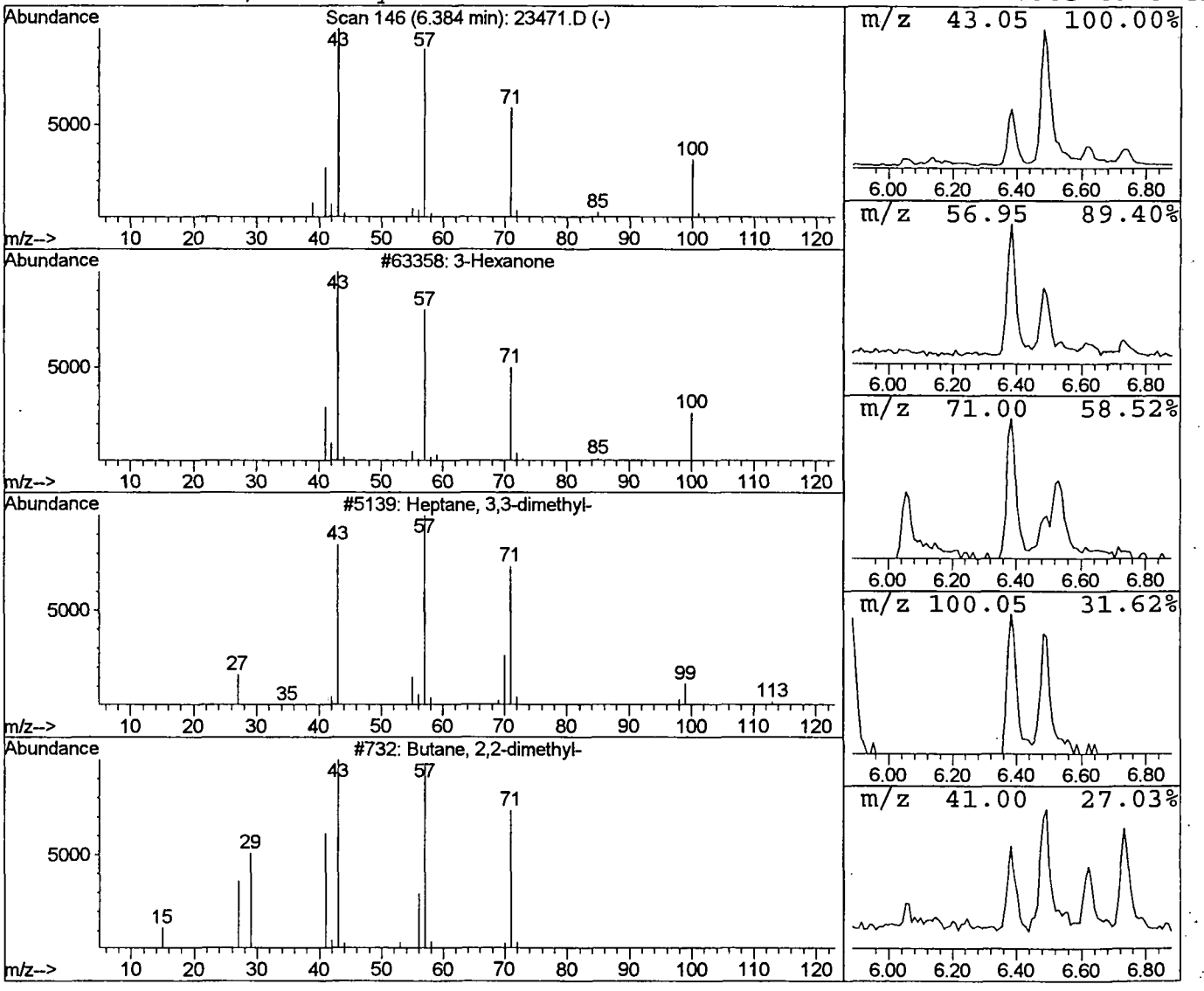
Quant Method : C:\HPCHEM\1\METHODS\NP091101.M (RTE Integrator)
Title : 209 625/TCLP calibration table 7/16/01
Library : C:\DATABASE\NBS75K.L

****

Peak Number 1 3-Hexanone Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.38	0.53 ug/l	51102	1,4-Dichlorobenzene-d4	11.76

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	3-Hexanone	100	C6H12O	000589-38-8	87
2	Heptane, 3,3-dimethyl-	128	C9H20	004032-86-4	59
3	Butane, 2,2-dimethyl-	86	C6H14	000075-83-2	53
4	3-Pentanone, 2-methyl-	100	C6H12O	000565-69-5	43



Library Search Compound Report

Data File : C:\HPCHEM\1\DATA\OCT0901\23471.D
Acq On : 10 Oct 2001 6:53 am
Sample : gc/ms unknown
Misc :
MS Integration Params: LSCINT.P

Vial: 19
Operator: 209 GALOS
Inst : GC/MS 209
Multiplr: 1.00

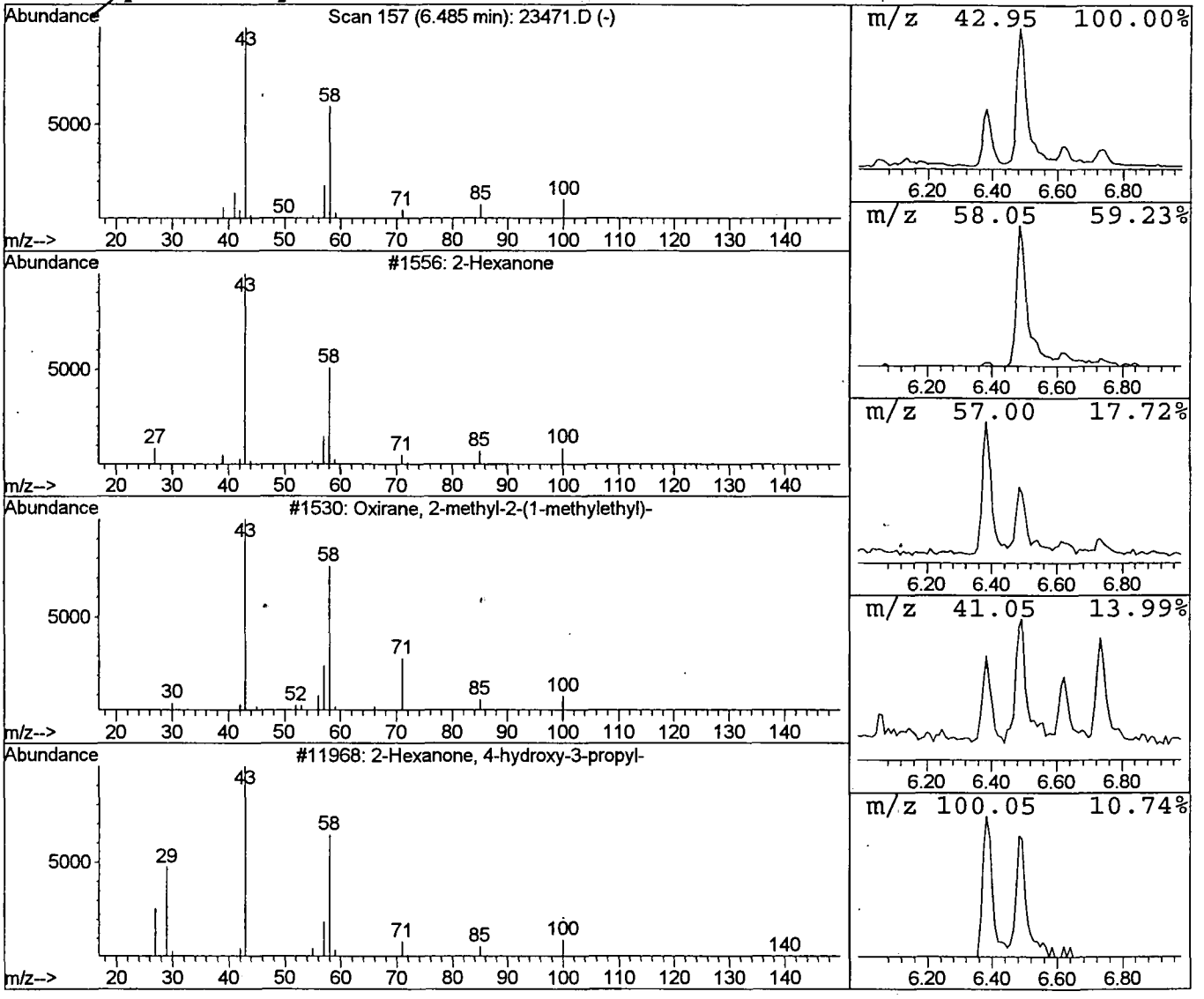
Quant Method : C:\HPCHEM\1\METHODS\NP091101.M (RTE Integrator)
Title : 209 625/TCLP calibration table 7/16/01
Library : C:\DATABASE\NBS75K.L

****

Peak Number 2 2-Hexanone Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.49	1.15 ug/l	110113	1,4-Dichlorobenzene-d4	11.76

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Hexanone	100	C6H12O	000591-78-6	91
2	Oxirane, 2-methyl-2-(1-methylethyl)	100	C6H12O	072221-03-5	72
3	2-Hexanone, 4-hydroxy-3-propyl-	158	C9H18O2	062338-17-4	64
4	Methyl Isobutyl Ketone	100	C6H12O	000108-10-1	52



Library Search Compound Report

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Acq On : 10 Oct 2001 6:53 am
Sample : gc/ms unknown
Misc :
MS Integration Params: LSCINT.P

Vial: 19
Operator: 209 GALOS
Inst : GC/MS 209
Multiplr: 1.00

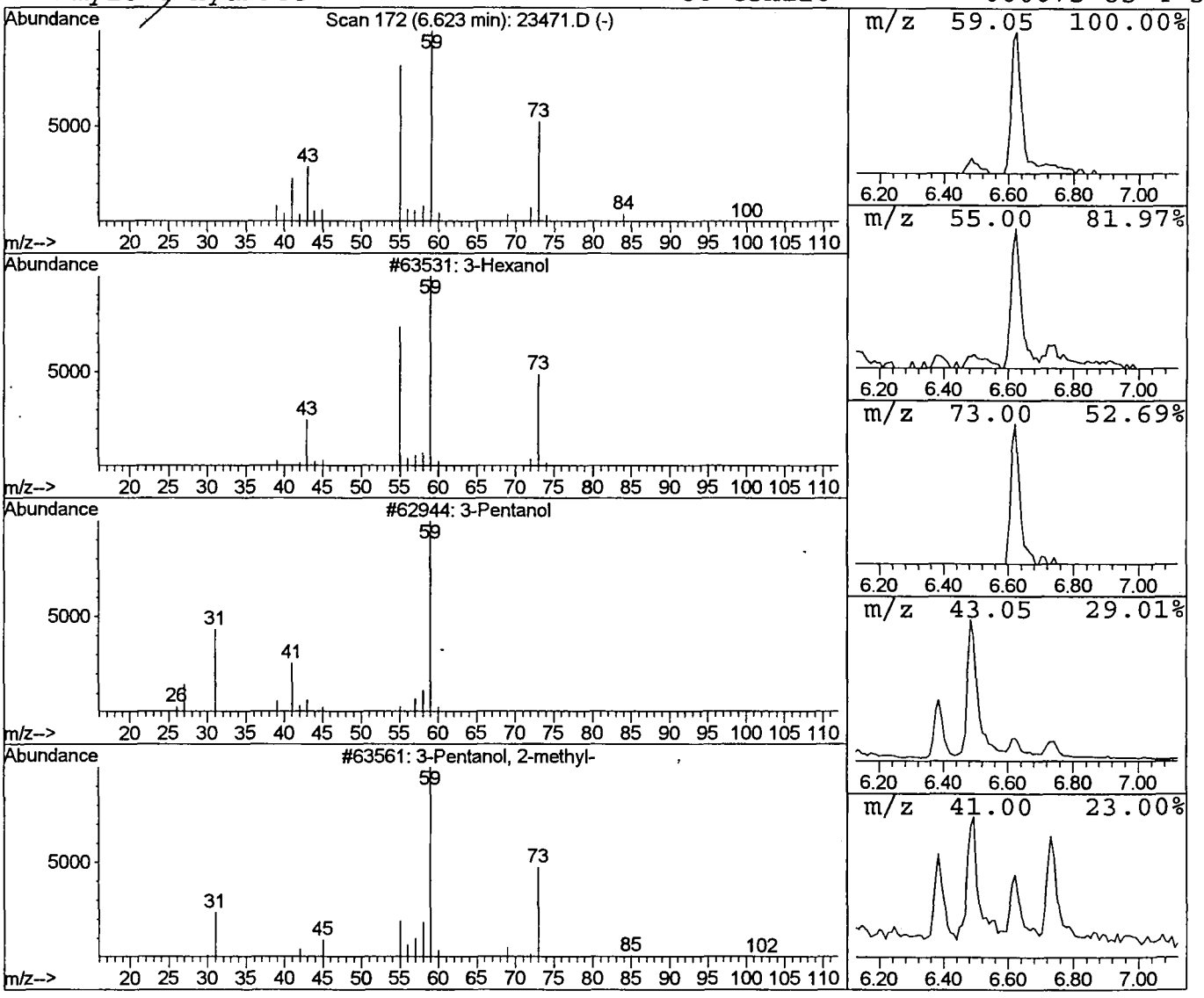
Quant Method : C:\HPCHEM\1\METHODS\NP091101.M (RTE Integrator)
Title : 209 625/TCLP calibration table 7/16/01
Library : C:\DATABASE\NBS75K.L

****

Peak Number 3 3-Hexanol — Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.62	0.51 ug/l	48469	1,4-Dichlorobenzene-d4	11.76

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	3-Hexanol	102	C6H14O	000623-37-0	90
2	3-Pentanol	88	C5H12O	000584-02-1	40
3	3-Pentanol, 2-methyl-	102	C6H14O	000565-67-3	39
4	Amylene Hydrate	88	C5H12O	000075-85-4	38



Library Search Compound Report

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Acq On : 10 Oct 2001 6:53 am
Sample : gc/ms unknown
Misc :
MS Integration Params: LSCINT.P

Vial: 19
Operator: 209 GALOS
Inst : GC/MS 209
Multiplr: 1.00

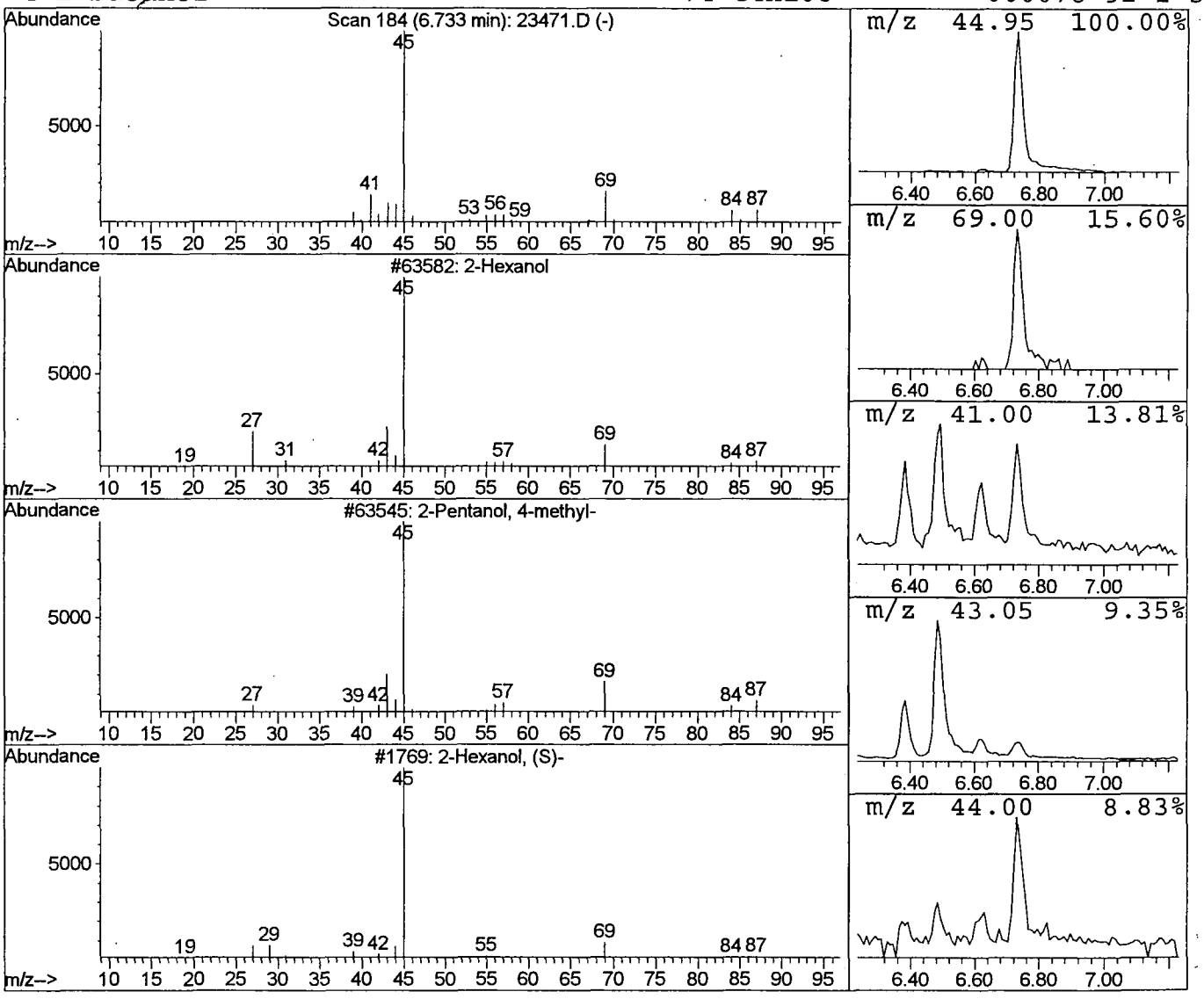
Quant Method : C:\HPCHEM\1\METHODS\NP091101.M (RTE Integrator)
Title : 209 625/TCLP calibration table 7/16/01
Library : C:\DATABASE\NBS75K.L

****

Peak Number 4 2-Hexanol — Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.73	0.68 ug/l	65171	1,4-Dichlorobenzene-d4	11.76

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	74
3	2-Hexanol, (S)-	102	C6H14O	052019-78-0	64
4	2-Butanol	74	C4H10O	000078-92-2	50



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Misc :
MS Integration Params: LSCINT.P

Vial: 19
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Inst : GC/MS 209
Multiplr: 1.00

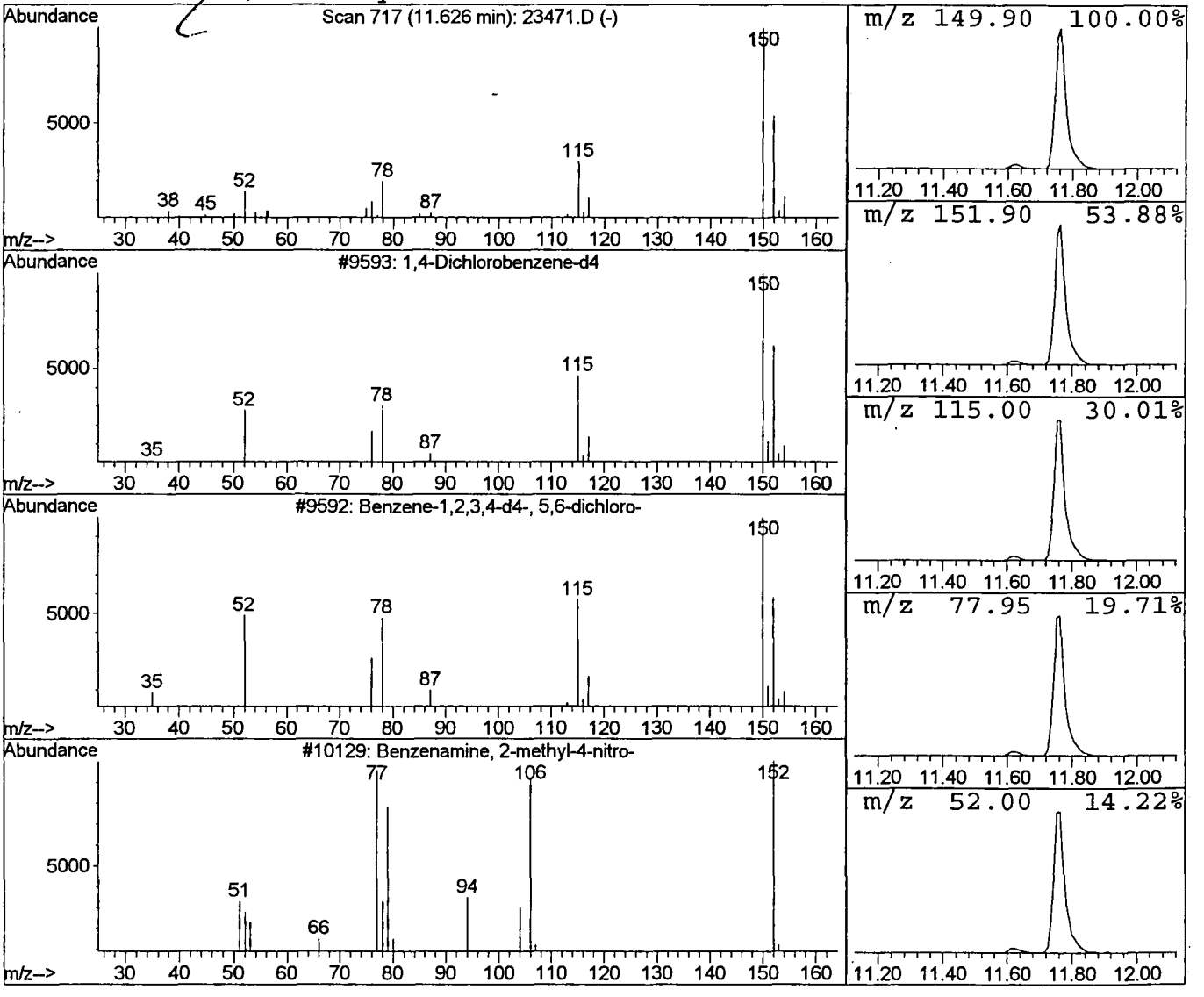
Quant Method : C:\HPCHEM\1\METHODS\NP091101.M (RTE Integrator)
Title : 209 625/TCLP calibration table 7/16/01
Library : C:\DATABASE\NBS75K.L

****

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.63	0.53 ug/l	50861	1,4-Dichlorobenzene-d4	11.76

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	33
3	Benzenamine, 2-methyl-4-nitro-	152	C7H8N2O2	000099-52-5	22
4	Benzenamine, 2-methyl-5-nitro-	152	C7H8N2O2	000099-55-8	14



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

Vial: 19
Operator: 209 GALOS
Inst : GC/MS 209
Multiplr: 1.00

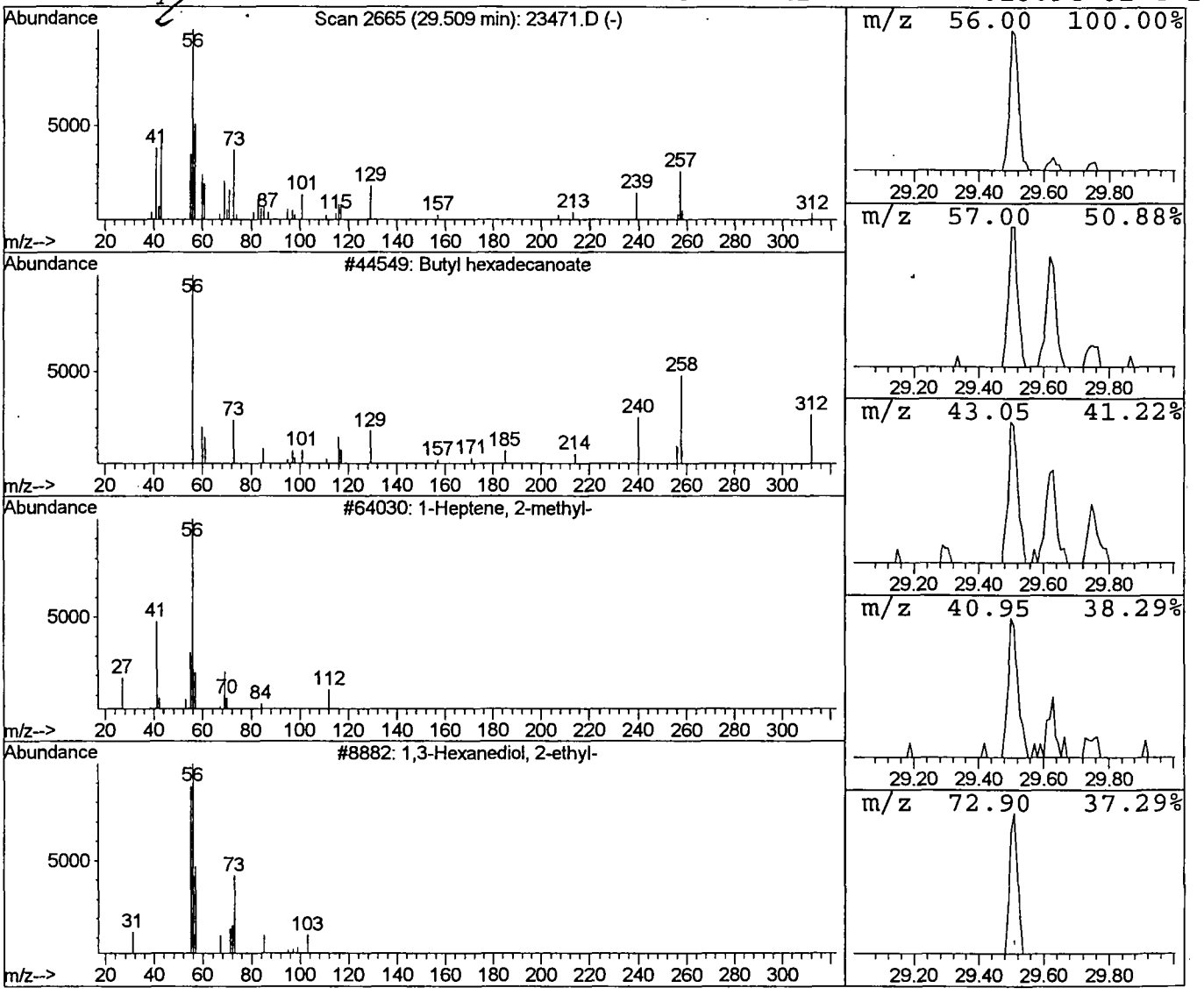
Quant Method : C:\HPCHEM\1\METHODS\NP091101.M (RTE Integrator)
Title : 209 625/TCLP calibration table 7/16/01
Library : C:\DATABASE\NBS75K.L

****

Peak Number 6 Butyl hexadecanoate Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
29.51	0.52 ug/l	44346	Chrysene-d12	33.12

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Butyl hexadecanoate	312	C20H40O2	000000-00-0	50
2	1-Heptene, 2-methyl-	112	C8H16	015870-10-7	27
3	1,3-Hexanediol, 2-ethyl-	146	C8H18O2	000094-96-2	23
4	2-Methyl-N-1-tridecene	196	C14H28	018094-01-4	16



Library Search Compound Report

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Misc :
MS Integration Params: LSCINT.P

Vial: 19
Operator: 209 GALOS
Inst : GC/MS 209
Multiplr: 1.00

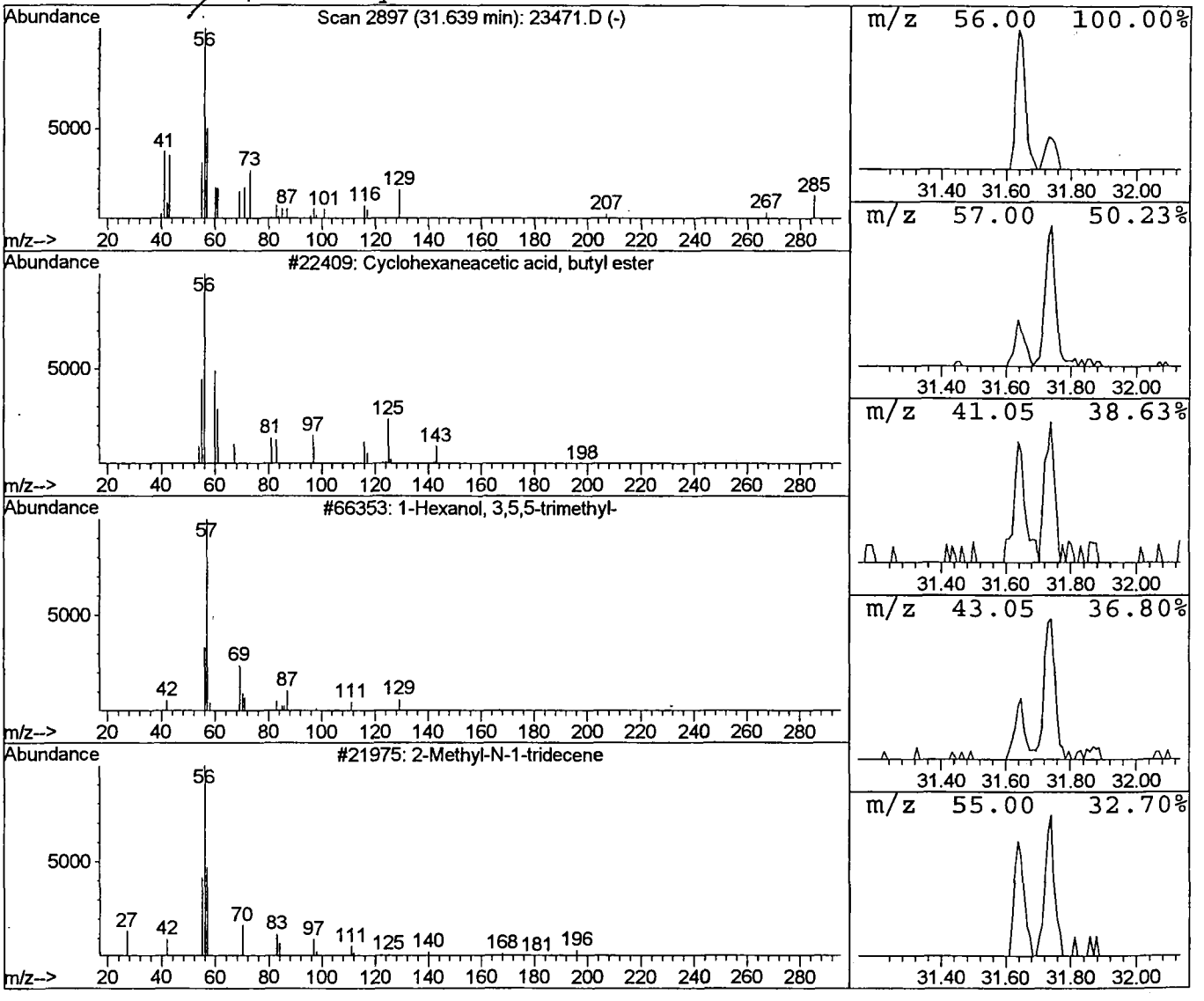
Quant Method : C:\HPCHEM\1\METHODS\NP091101.M (RTE Integrator)
Title : 209 625/TCLP calibration table 7/16/01
Library : C:\DATABASE\NBS75K.L

****

Peak Number 7 Cyclohexaneacetic acid, butyl Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
31.64	0.41 ug/l	35046	Chrysene-d12	33.12

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclohexaneacetic acid, butyl ester	198	C12H22O2	027948-12-5	38
2	1-Hexanol, 3,5,5-trimethyl-	144	C9H20O	003452-97-9	17
3	2-Methyl-N-1-tridecene	196	C14H28	018094-01-4	16
4	Pentadecane, 8-methylene-	224	C16H32	055668-09-2	16



Library Search Compound Report

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Acq On : 10 Oct 2001 6:53 am
Sample : gc/ms unknown
Misc :
MS Integration Params: LSCINT.P

Vial: 19
Operator: 209 GALOS
Inst : GC/MS 209
Multiplr: 1.00

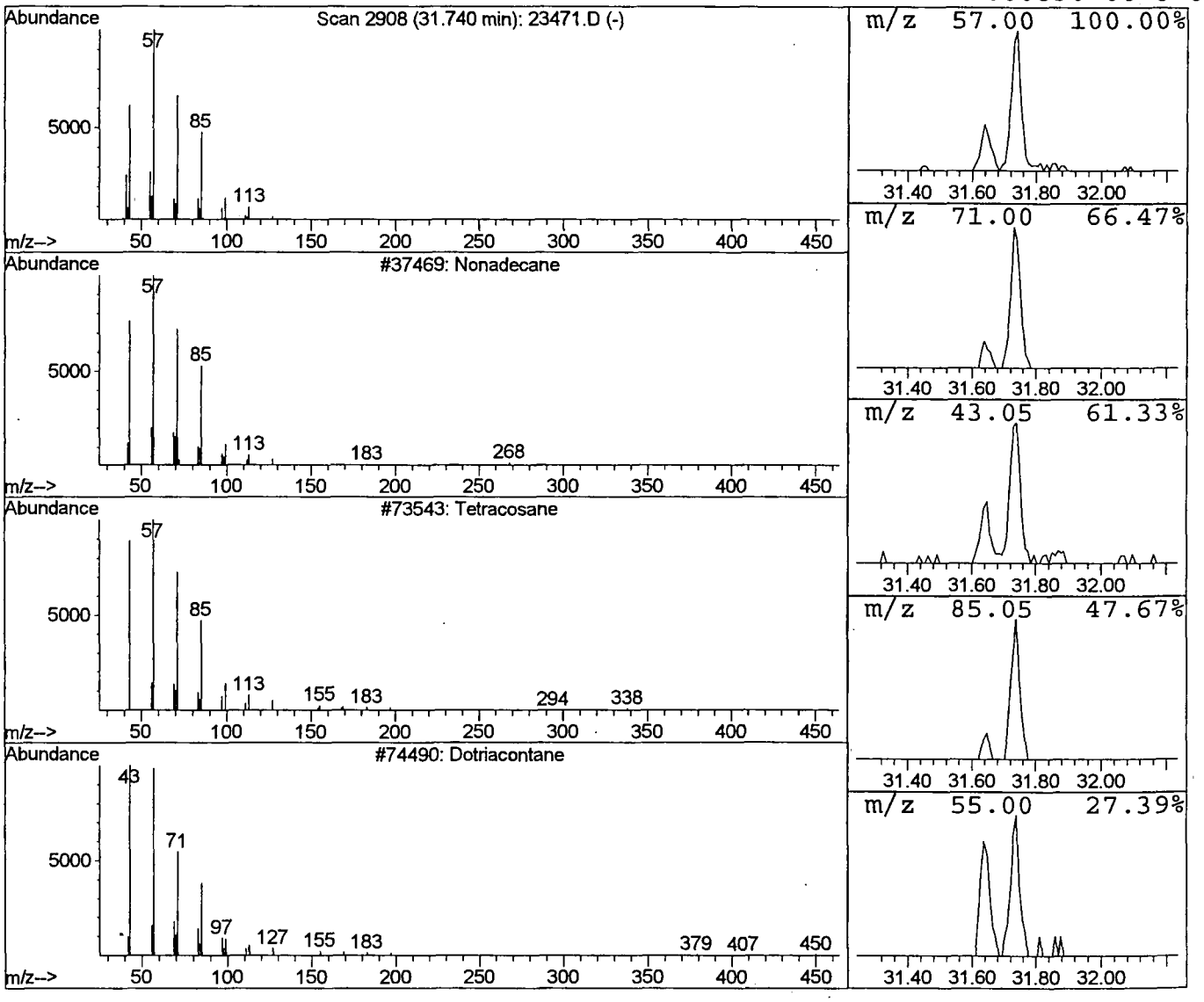
Quant Method : C:\HPCHEM\1\METHODS\NP091101.M (RTE Integrator)
Title : 209 625/TCLP calibration table 7/16/01
Library : C:\DATABASE\NBS75K.L

****

Peak Number 8 Nonadecane — Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
31.74	0.52 ug/l	44576	Chrysene-d12	33.12

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Nonadecane	268	C19H40	000629-92-5	91
2	Tetracosane	338	C24H50	000646-31-1	90
3	Dotriacontane	451	C32H66	000544-85-4	90
4	Hexatriacontane	507	C36H74	000630-06-8	86



Library Search Compound Report

Data File : C:\HPCHEM\1\DATA\OCT0901\23471.D
Acq On : 10 Oct 2001 6:53 am
Sample : gc/ms unknown
Misc :
MS Integration Params: LSCINT.P

Vial: 19
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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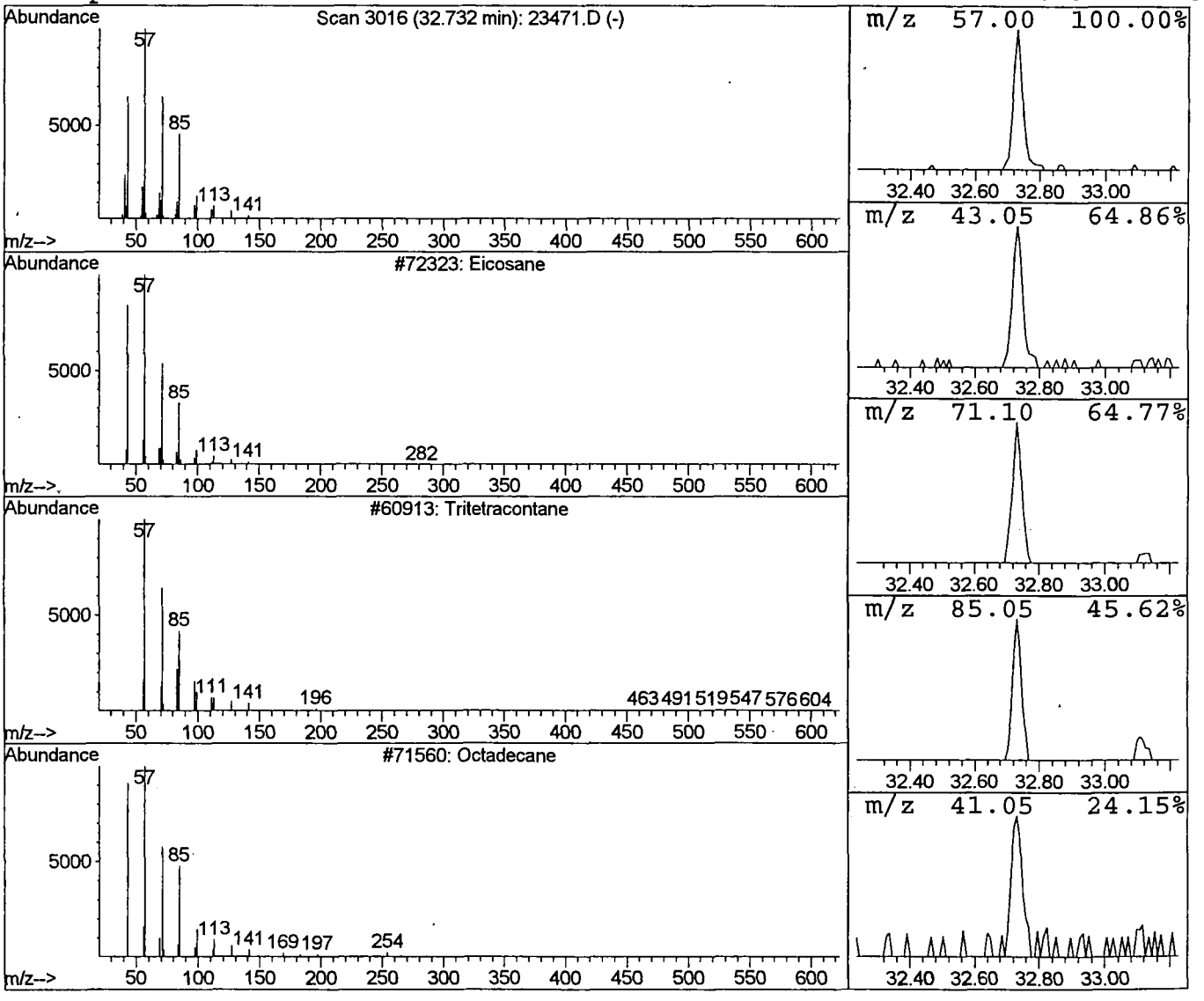
Peak Number 9 Eicosane — Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
32.73	0.52 ug/l	44327	Chrysene-d12	33.12

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Eicosane	282	C20H42	000112-95-8	91
2	Tritetracontane	605	C43H88	007098-21-7	90
3	Octadecane	254	C18H38	000593-45-3	90
4	Heptadecane	240	C17H36	000629-78-7	90



Library Search Compound Report

Data File : C:\HPCHEM\1\DATA\OCT0901\23471.D
Acq On : 10 Oct 2001 6:53 am
Sample : gc/ms unknown
Misc :
MS Integration Params: LSCINT.P

Vial: 19
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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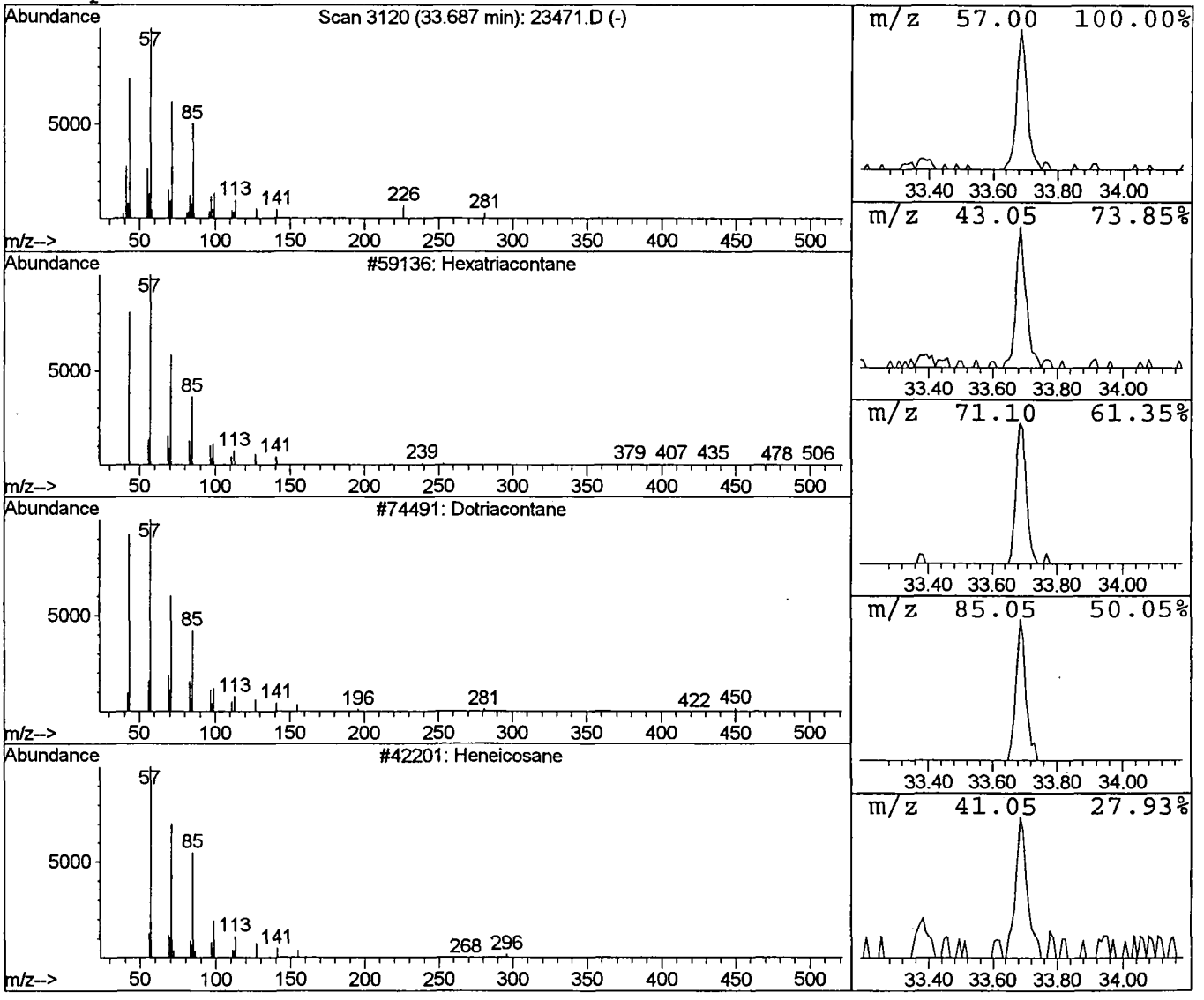
Peak Number 10 Hexatriacontane Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
33.69	0.64 ug/l	55225	Chrysene-d12	33.12

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Hexatriacontane	507	C36H74	000630-06-8	90
2	Dotriacontane	451	C32H66	000544-85-4	90
3	Heneicosane	296	C21H44	000629-94-7	90
4	Heptacosane	380	C27H56	000593-49-7	87



Library Search Compound Report

Data File : C:\HPCHEM\1\DATA\OCT0901\23471.D
Acq On : 10 Oct 2001 6:53 am
Sample : gc/ms unknown
Misc :
MS Integration Params: LSCINT.P

Vial: 19
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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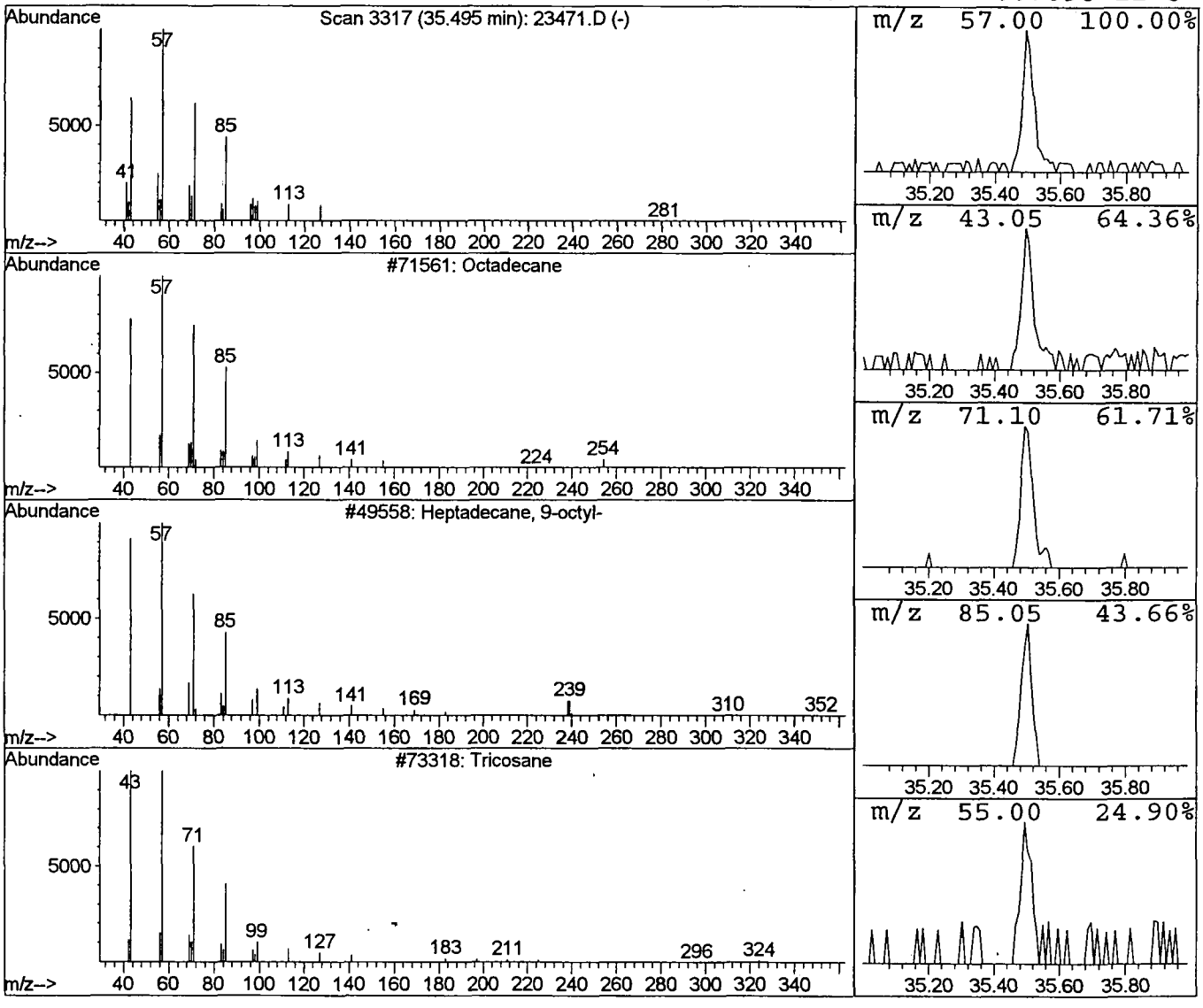
Peak Number 11 Octadecane — Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
35.50	0.41 ug/l	28532	Perylene-d12	37.23

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Octadecane	254	C18H38	000593-45-3	86
2	Heptadecane, 9-octyl-	352	C25H52	007225-64-1	86
3	Tricosane	324	C23H48	000638-67-5	83
4	Tetratetracontane	619	C44H90	007098-22-8	83



Tentatively Identified Compound (LSC) summary

Operator ID: 209 GALOS Date Acquired: 10 Oct 2001 6:53 am
Data File: C:\HPCHEM\1\DATA\OCT0901\23471.D
Name: gc/ms unknown
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	0.5	ug/l	51102	ISTD01	11.76	1913980	20.0
2-Hexanone	6.49	1.2	ug/l	110113	ISTD01	11.76	1913980	20.0
3-Hexanol	6.62	0.5	ug/l	48469	ISTD01	11.76	1913980	20.0
2-Hexanol	6.73	0.7	ug/l	65171	ISTD01	11.76	1913980	20.0
1,4-Dichlorobenzene	11.63	0.5	ug/l	50861	ISTD01	11.76	1913980	20.0
Butyl hexadecanoate	29.51	0.5	ug/l	44346	ISTD05	33.12	1713810	20.0
Cyclohexaneacetic ac	31.64	0.4	ug/l	35046	ISTD05	33.12	1713810	20.0
Nonadecane	31.74	0.5	ug/l	44576	ISTD05	33.12	1713810	20.0
Eicosane	32.73	0.5	ug/l	44327	ISTD05	33.12	1713810	20.0
Hexatriacontane	33.69	0.6	ug/l	55225	ISTD05	33.12	1713810	20.0
Octadecane	35.50	0.4	ug/l	28532	ISTD06	37.23	1393950	20.0

23471.D NP091101.M Thu Oct 11 13:15:41 2001

Comments for Sample AD23471 - Analysis $GCMS

Westchester Dept. of Labs & Research
User: Vinci, David L
Date: 10-19-2001 Time: 17:11:38

A scan of the extract prepared for the 508 analysis, from 35 - 550 amu using a mass spectrometer, reveals the presence of non-target compounds. However, these compounds were observed in previous Laboratory Blanks, though not in the Laboratory Blank associated with this sample.