User: Galos, Marie
Westchester Dept. of Labs & Research
Date: 10/11/2001 Time: 08:24:19

Sample: AD22653
Analysis: $GCMS GC/MS Scan For Unknowns
Started: 10/11/01 08:23 Ended: 10/11/01 08:23 Analyst: MG
Page: 1

	Component Name	Viol	Result
1	Other unknown compounds		see comment

Quantitation Report (QT Reviewed)

Data File : C:\HPCHEM\1\DATA\SEP2801\22653.D Vial: 12
Acq On : 28 Sep 2001 11:13 pm Operator: 209 GALOS
Sample : gc/ms unknown Inst : GC/MS 209
Misc : Multiplr: 1.00
MS Integration Params: RTEINT.P
Quant Time: Oct 3 14:27 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	11.79	152	483889	20.00	ug/l	-0.12
19) Naphthalene-d8	15.41	136	1879031m	20.00	ug/l	-0.12
31) Acenaphthene-d10	20.70	164	891776	20.00	ug/l	-0.12
49) Phenanthrene-d10	25.12	188	1277714	20.00	ug/l	-0.12
59) Chrysene-d12	33.16	240	959728	20.00	ug/l	-0.12
68) Perylene-d12	37.28	264	735700	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	105	0.00	ug/l	0.37
Spiked Amount 75.000			Recovery =	0.00%		
7) 1,2-Dichlorobenzene-d4	12.30	152	5318	0.23	ug/l	-0.13
Spiked Amount 50.000			Recovery =	0.46%		
20) Nitrobenzene-d5	0.00	82	0	0.00	ug/l	
Spiked Amount 50.000			Recovery =	0.00%		
32) 2-Fluorobiphenyl	18.82	172	1930	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	5.18	79	302	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

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Last Update : Thu Sep 13 12:47:04 2001
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DataAcq Meth : NP091101

	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	22.18	149	386	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	25.12	178	408	N.D.	
56)	Anthracene	25.12	178	408	N.D.	
57)	Di-n-butylphthalate	27.11	149	3788	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	33.16	228	2260	N.D.	
66)	Bis(2-ethylhexyl)phthalate	33.42	149	11379	N.D.	
67)	Chrysene	33.16	228	2260	N.D.	
69)	Di-n-Octylphthalate	35.16	149	2460	N.D.	
70)	Benzo(b)fluoranthene	36.19	252	117	N.D.	

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Misc : Multiplr: 1.00
MS Integration Params: RTEINT.P
Quant Time: Oct 3 14:27 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

	Compound	R.T.	QIon	Response	Conc Unit	Qvalue
71)	Benzo(k)fluoranthene	36.25	252	193	N.D.	
72)	Benzo(a)pyrene	37.28	252	2698	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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Sample : gc/ms unknown
Misc :
MS Integration Params: RTEINT.P
Quant Time: Oct 3 14:27 2001

Vial: 12
Operator: 209 GALOS
Inst : GC/MS 209
Multiplr: 1.00
Quant Results File: NP091101.RES

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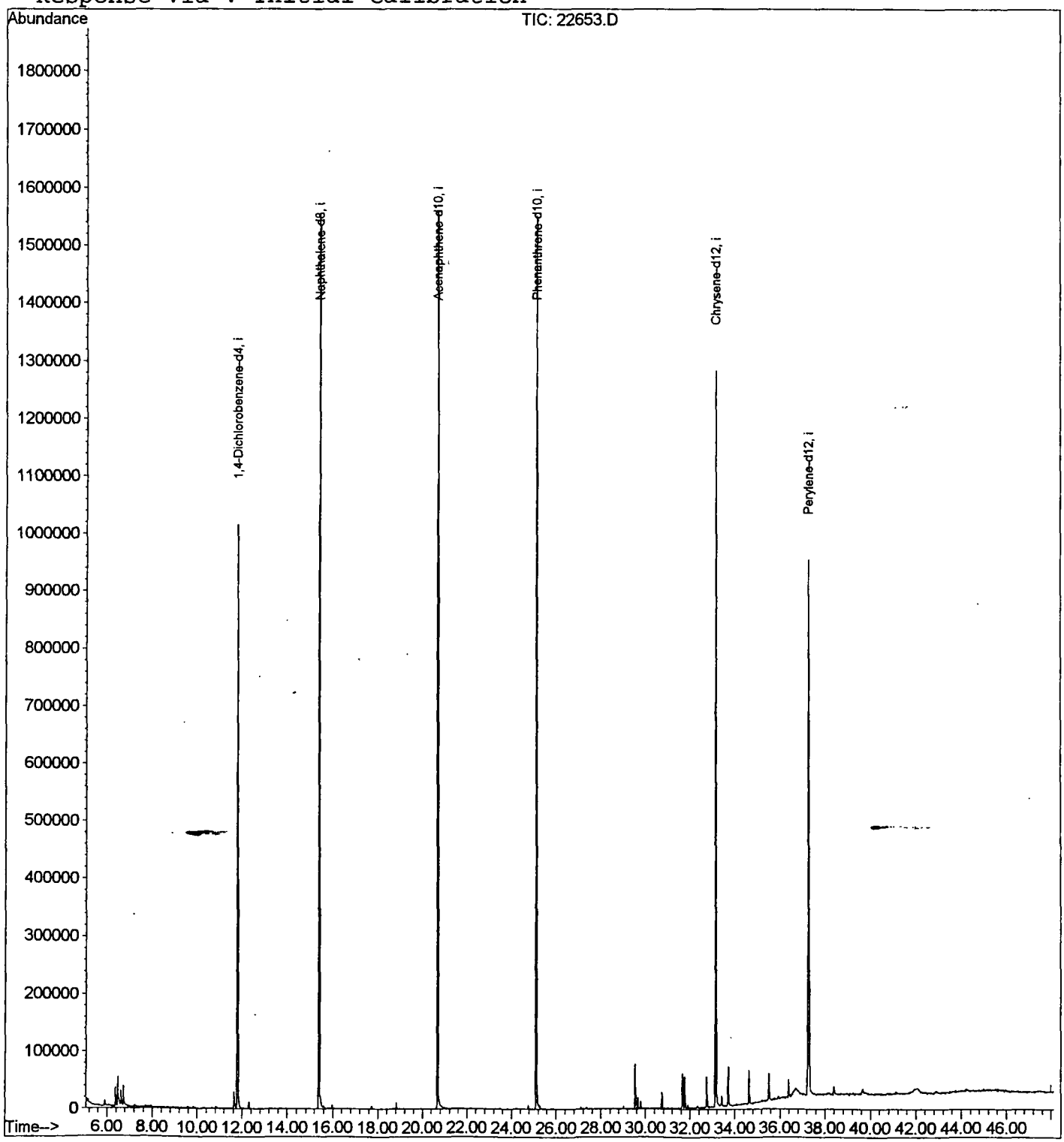
Title : 209 625/TCLP calibration table 7/16/01
Last Update : Thu Sep 13 12:47:04 2001
Response via : Initial Calibration



LSC Area Percent Report

Data File : C:\HPCHEM\1\DATA\SEP2801\22653.D Vial: 12
Acq On : 28 Sep 2001 11:13 pm 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

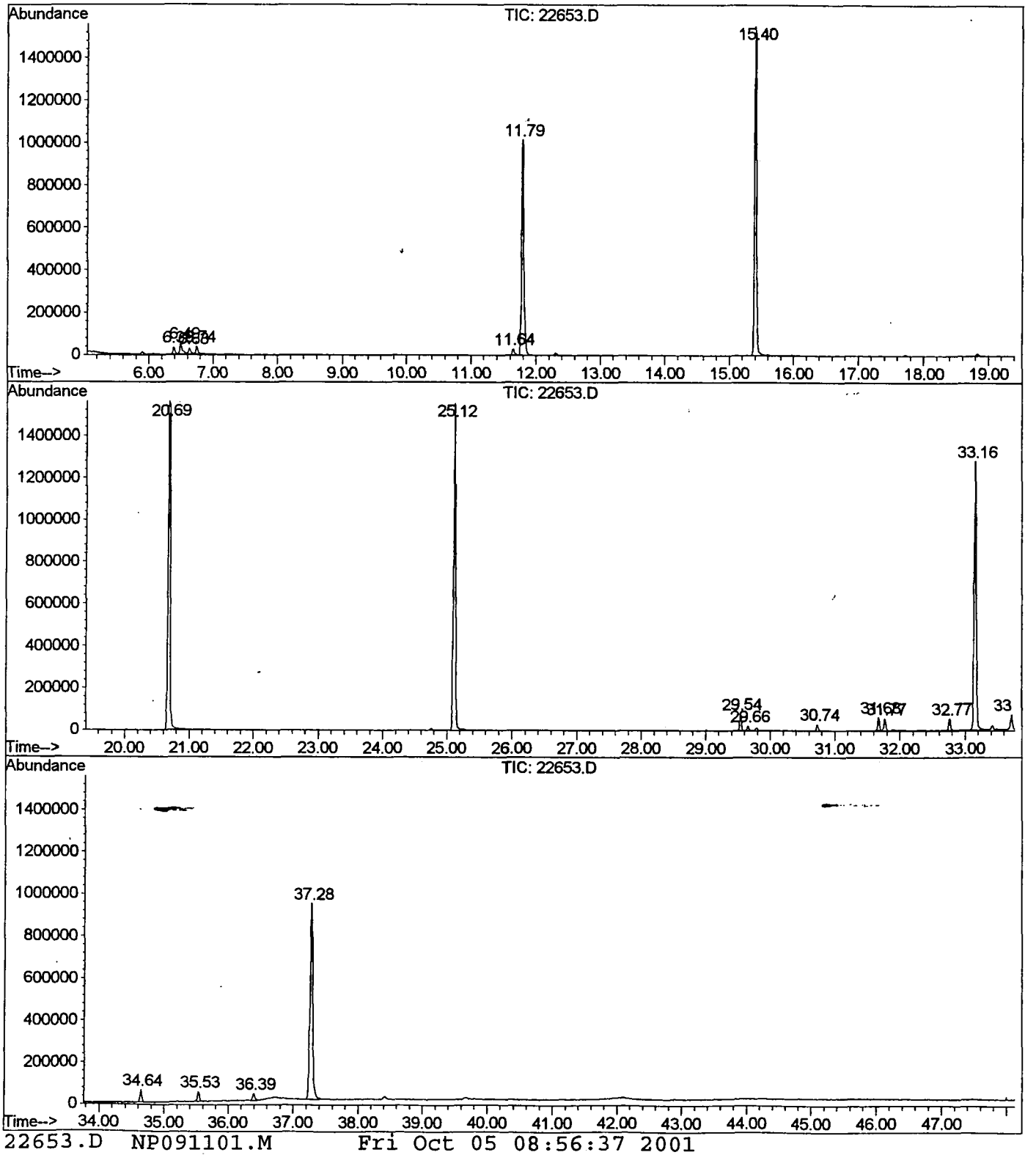
peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	peak area	peak % max.	% of total
1	6.392	142	147	153	rBV	32891	69191	1.96%	0.347%
2	6.493	154	158	169	rVV2	51730	133042	3.76%	0.667%
3	6.631	169	173	178	rVV	26838	56478	1.60%	0.283%
4	6.741	181	185	196	rVB	34729	74340	2.10%	0.373%
5	11.644	714	719	728	rBV	28344	67113	1.90%	0.336%
6	11.790	729	735	749	rVB	1014262	2500964	70.70%	12.532%
7	15.398	1122	1128	1145	rBV	1551320	3450459	97.54%	17.290%
8	20.686	1697	1704	1717	rBV	1559467	3537645	100.00%	17.727%
9	25.120	2180	2187	2199	rBV	1549759	3316878	93.76%	16.620%
10	29.536	2663	2668	2675	rBV	78794	150248	4.25%	0.753%
11	29.655	2675	2681	2686	rVV2	19803	37917	1.07%	0.190%
12	30.739	2793	2799	2805	rBV	28180	57249	1.62%	0.287%
13	31.675	2897	2901	2905	rVB2	60160	112947	3.19%	0.566%
14	31.767	2907	2911	2917	rBV	53950	109618	3.10%	0.549%
15	32.767	3013	3020	3027	rBV	53799	116242	3.29%	0.582%
16	33.162	3056	3063	3076	rBV2	1278817	2976965	84.15%	14.917%
17	33.722	3114	3124	3130	rBV2	69182	160266	4.53%	0.803%
18	34.640	3220	3224	3230	rVB	57497	111676	3.16%	0.560%
19	35.531	3316	3321	3326	rBV	46516	101310	2.86%	0.508%
20	36.394	3410	3415	3423	rBV	31498	80479	2.27%	0.403%
21	37.284	3503	3512	3531	rBV2	929018	2735723	77.33%	13.708%

Sum of corrected areas: 19956750

22653.D NP091101.M Fri Oct 05 08:56:35 2001

LSC Report - Integrated Chromatogram

File : C:\HPCHEM\1\DATA\SEP2801\22653.D
Operator : 209 GALOS
Acquired : 28 Sep 2001 11:13 pm using AcqMethod NP091101
Instrument : GC/MS 209
Sample Name: gc/ms unknown
Misc Info :
Vial Number: 12
Quant File :NP091101.RES (RTE Integrator)



22653.D NP091101.M Fri Oct 05 08:56:37 2001

Library Search Compound Report

Data File : C:\HPCHEM\1\DATA\SEP2801\22653.D
Acq On : 28 Sep 2001 11:13 pm
Sample : gc/ms unknown
Misc :
MS Integration Params: LSCINT.P

Vial: 12
Operator: 209 GALOS
Inst : GC/MS 209
Multiplr: 1.00

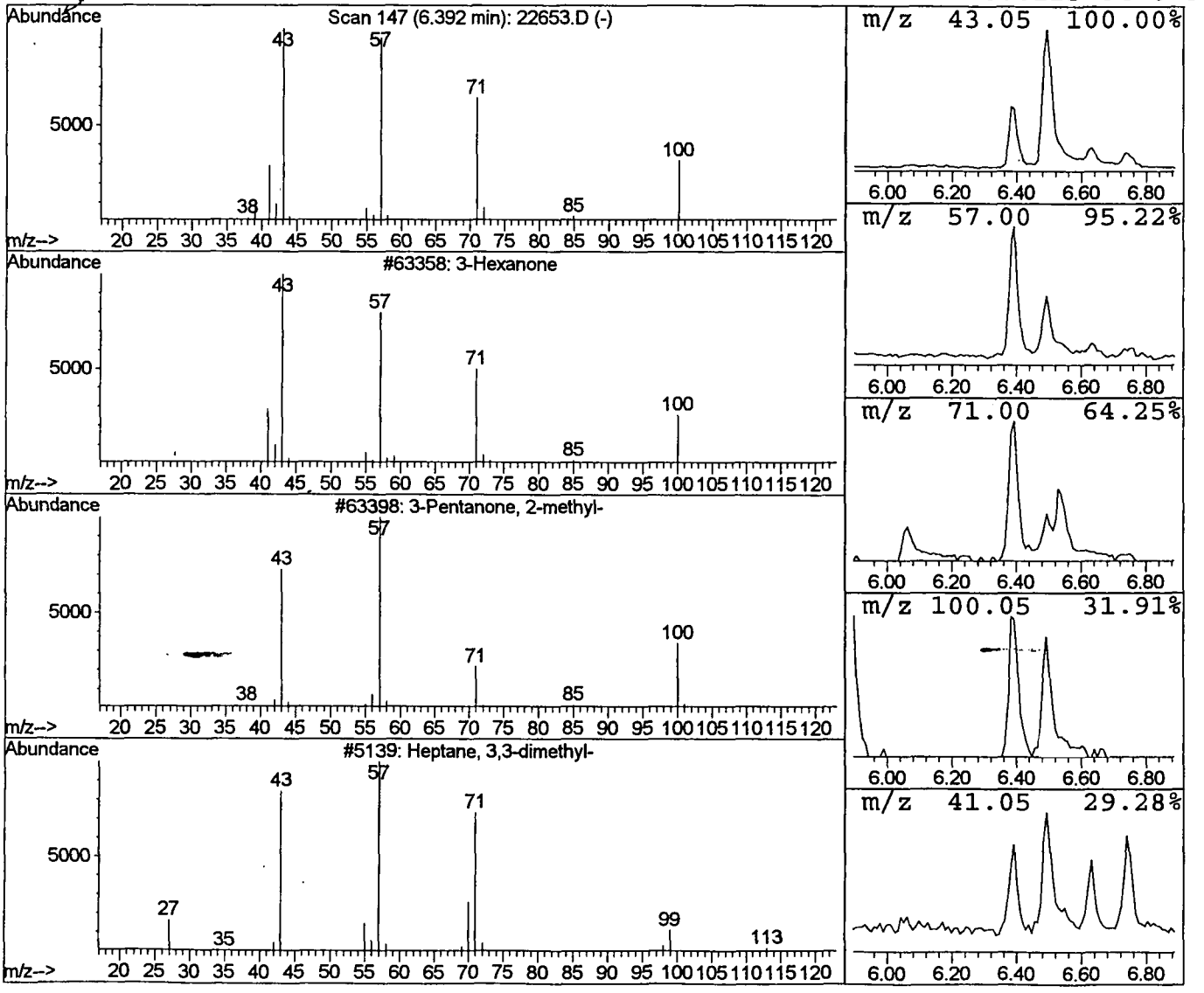
Quant Method : C:\HPCHEM\1\METHODS\NP091101.M (RTE Integrator)
Title : 209 625/TCLP calibration table 7/16/01
Library : C:\DATABASE\NBS75K.L

****

Peak Number 1 3-Hexanone Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.39	0.55 ug/l	69191	1,4-Dichlorobenzene-d4	11.79

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	37
3	Heptane, 3,3-dimethyl-	128	C9H20	004032-86-4	36
4	2,4-Pentanedione	100	C5H8O2	000123-54-6	35



Library Search Compound Report

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Acq On : 28 Sep 2001 11:13 pm
Sample : gc/ms unknown
Misc :
MS Integration Params: LSCINT.P

Vial: 12
Operator: 209 GALOS
Inst : GC/MS 209
Multiplr: 1.00

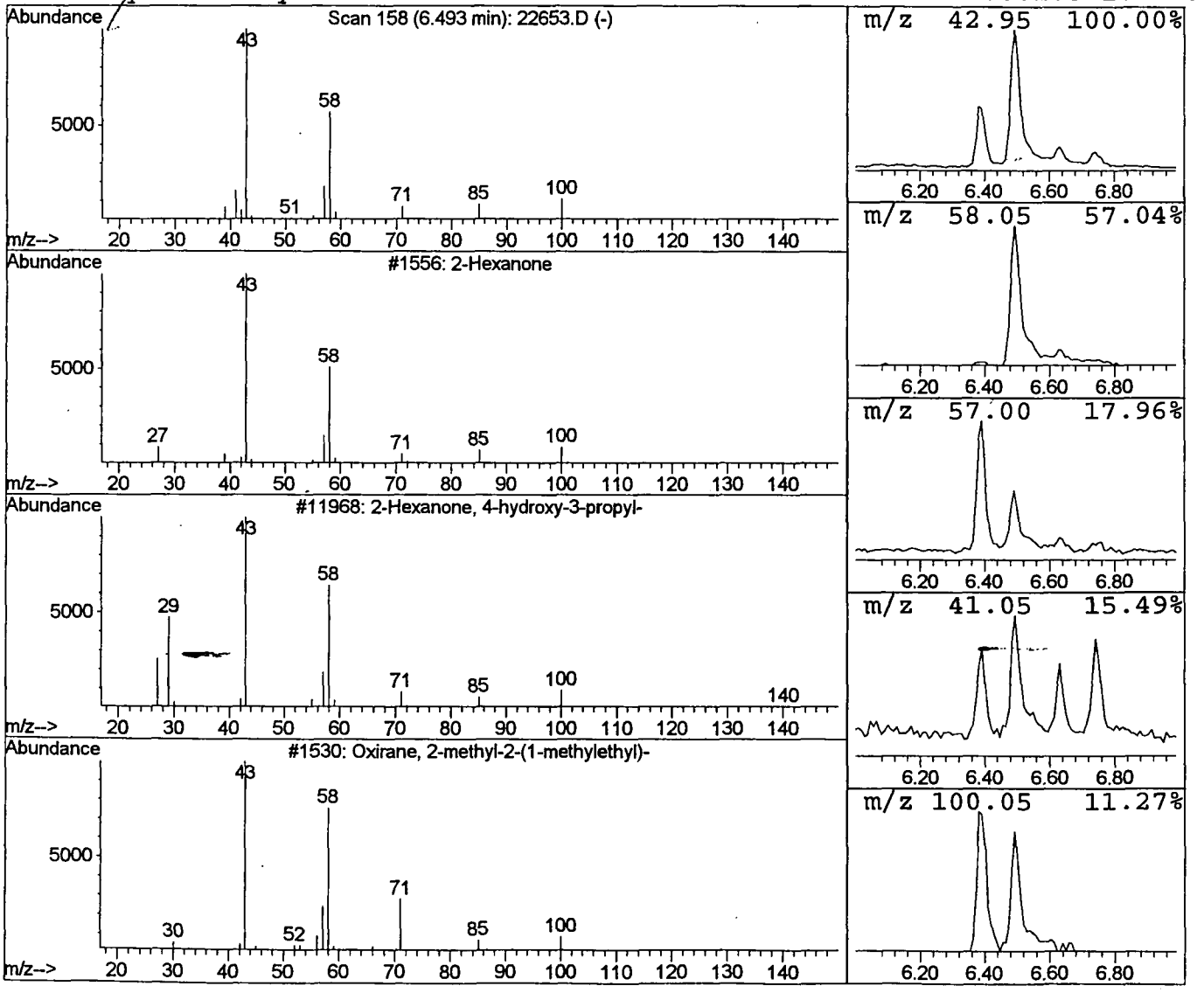
Quant Method : C:\HPCHEM\1\METHODS\NP091101.M (RTE Integrator)
Title : 209 625/TCLP calibration table 7/16/01
Library : C:\DATABASE\NBS75K.L

****

Peak Number 2 2-Hexanone Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.49	1.06 ug/l	133042	1,4-Dichlorobenzene-d4	11.79

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Hexanone	100	C6H12O	000591-78-6	91
2	2-Hexanone, 4-hydroxy-3-propyl-	158	C9H18O2	062338-17-4	83
3	Oxirane, 2-methyl-2-(1-methylethyl)	100	C6H12O	072221-03-5	78
4	Methyl Isobutyl Ketone	100	C6H12O	000108-10-1	58



Library Search Compound Report

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Acq On : 28 Sep 2001 11:13 pm
Sample : gc/ms unknown
Misc :
MS Integration Params: LSCINT.P

Vial: 12
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
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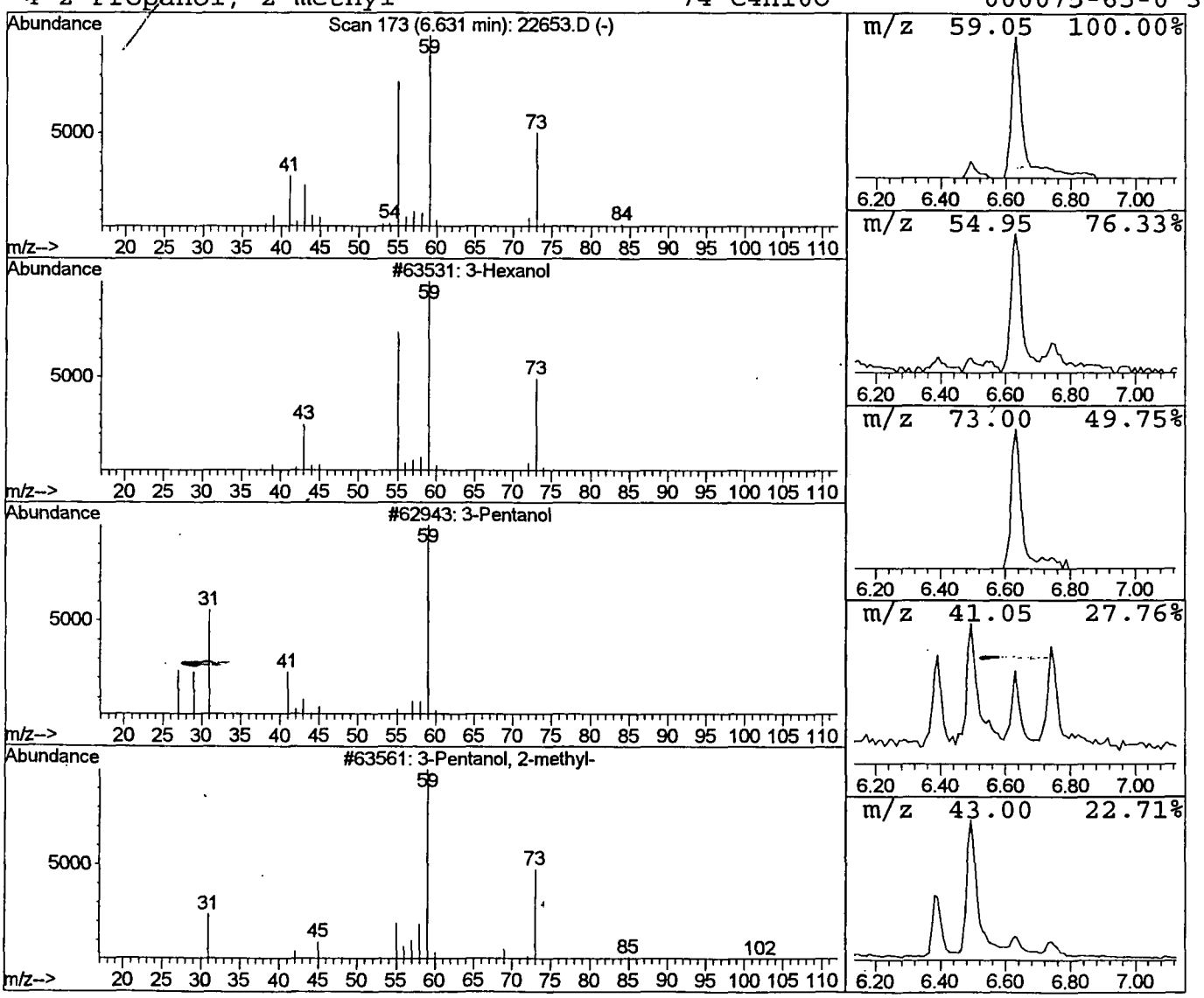
****

Peak Number 3 3-Hexanol Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.63	0.45 ug/l	56478	1,4-Dichlorobenzene-d4	11.79

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	45
3	3-Pentanol, 2-methyl-	102	C6H14O	000565-67-3	38
4	2-Propanol, 2-methyl-	74	C4H10O	000075-65-0	36



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Acq On : 28 Sep 2001 11:13 pm
Sample : gc/ms unknown
Misc :
MS Integration Params: LSCINT.P

Vial: 12
Operator: 209 GALOS
Inst : GC/MS 209
Multiplr: 1.00

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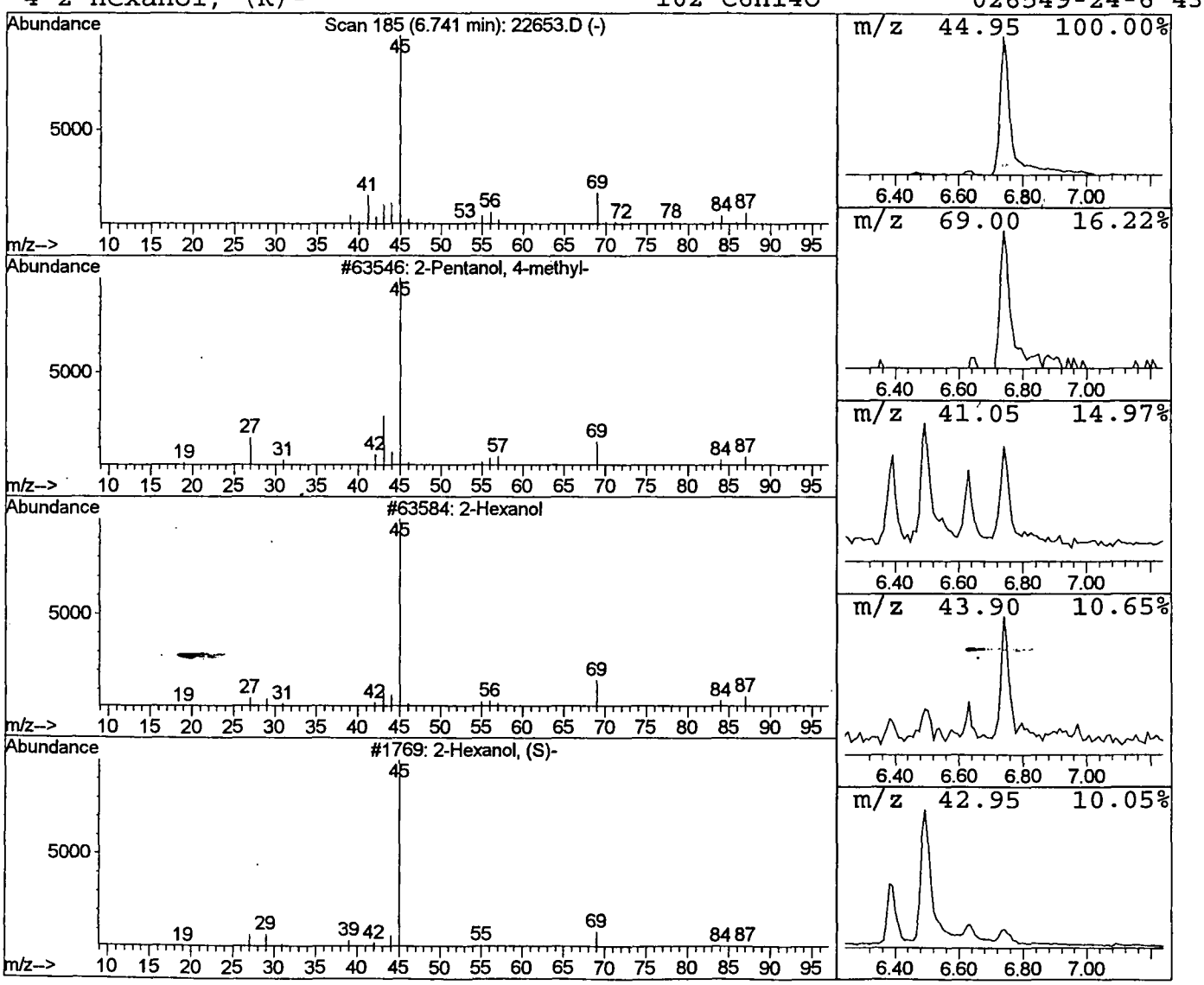
****

Peak Number 4 2-Pentanol, 4-methyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.74	0.59 ug/l	74340	1,4-Dichlorobenzene-d4	11.79

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Pentanol, 4-methyl-	102	C6H14O	000108-11-2	83
2	2-Hexanol	102	C6H14O	000626-93-7	83
3	2-Hexanol, (S)-	102	C6H14O	052019-78-0	64
4	2-Hexanol, (R)-	102	C6H14O	026549-24-6	43



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

Vial: 12
Operator: 209 GALOS
Inst : GC/MS 209
Multiplr: 1.00

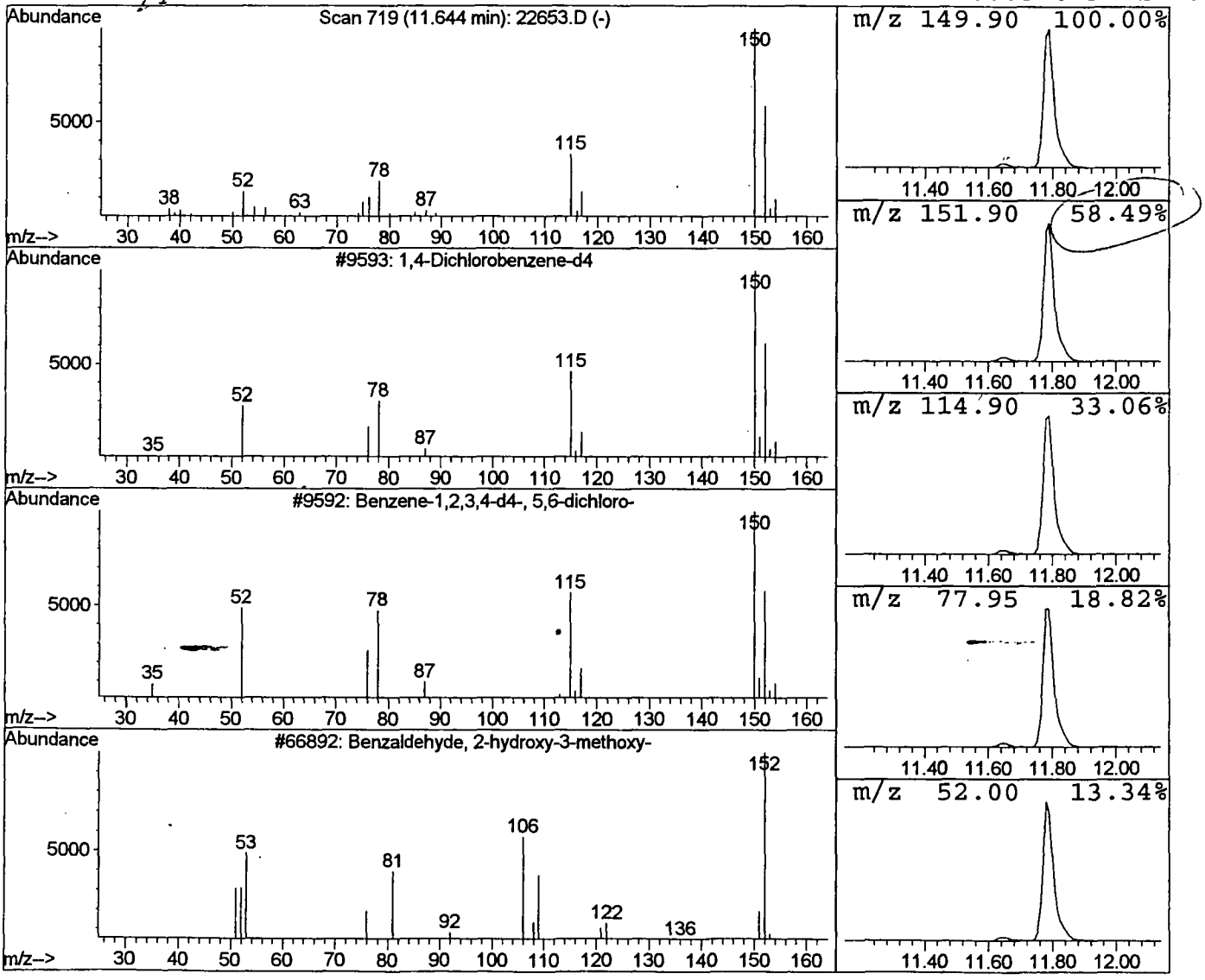
Quant Method : C:\HPCHEM\1\METHODS\NP091101.M (RTE Integrator)
Title : 209 625/TCLP calibration table 7/16/01
Library : C:\DATABASE\NBS75K.L

****

Peak Number 5 1,4-Dichlorobenzene-d4 Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.64	0.54 ug/l	67113	1,4-Dichlorobenzene-d4	11.79

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,4-Dichlorobenzene-d4	150	C6Cl2D4	003855-82-1	53
2	Benzene-1,2,3,4-d4-, 5,6-dichloro-	150	C6Cl2D4	002199-69-1	33
3	Benzaldehyde, 2-hydroxy-3-methoxy-	152	C8H8O3	000148-53-8	10
4	2-Methyl-6-nitroaniline	152	C7H8N2O2	000570-24-1	10



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Misc :
MS Integration Params: LSCINT.P

Vial: 12
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Multiplr: 1.00

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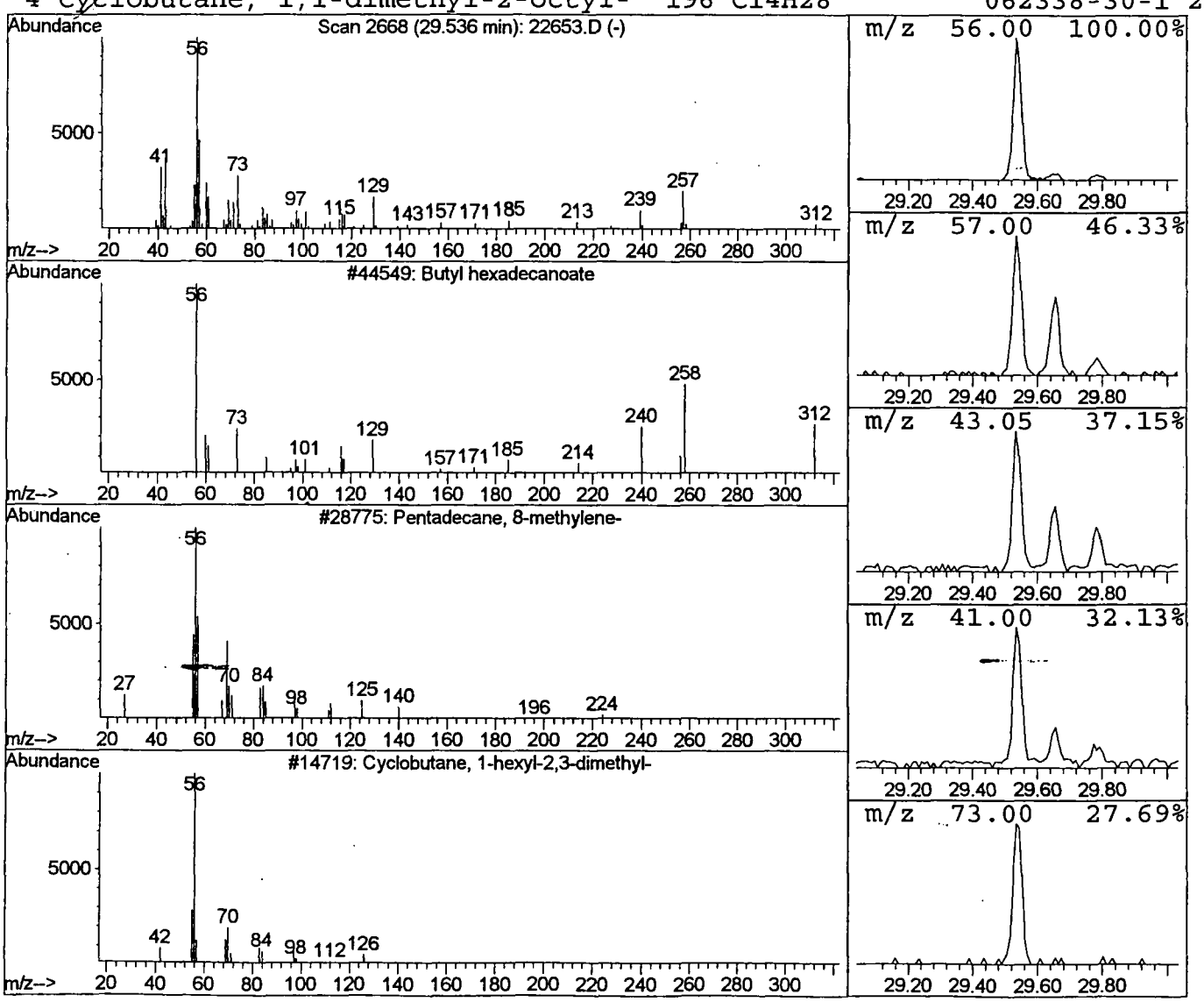
Library : C:\DATABASE\NBS75K.L

****

Peak Number 6 Butyl hexadecanoate Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
29.54	1.01 ug/l	150248	Chrysene-d12	33.16

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Butyl hexadecanoate	312	C20H40O2	000000-00-0	64
2	Pentadecane, 8-methylene-	224	C16H32	055668-09-2	25
3	Cyclobutane, 1-hexyl-2,3-dimethyl-	168	C12H24	055170-84-8	25
4	Cyclobutane, 1,1-dimethyl-2-octyl-	196	C14H28	062338-30-1	25



Library Search Compound Report

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Acq On : 28 Sep 2001 11:13 pm
Sample : gc/ms unknown
Misc :
MS Integration Params: LSCINT.P

Vial: 12
Operator: 209 GALOS
Inst : GC/MS 209
Multiplr: 1.00

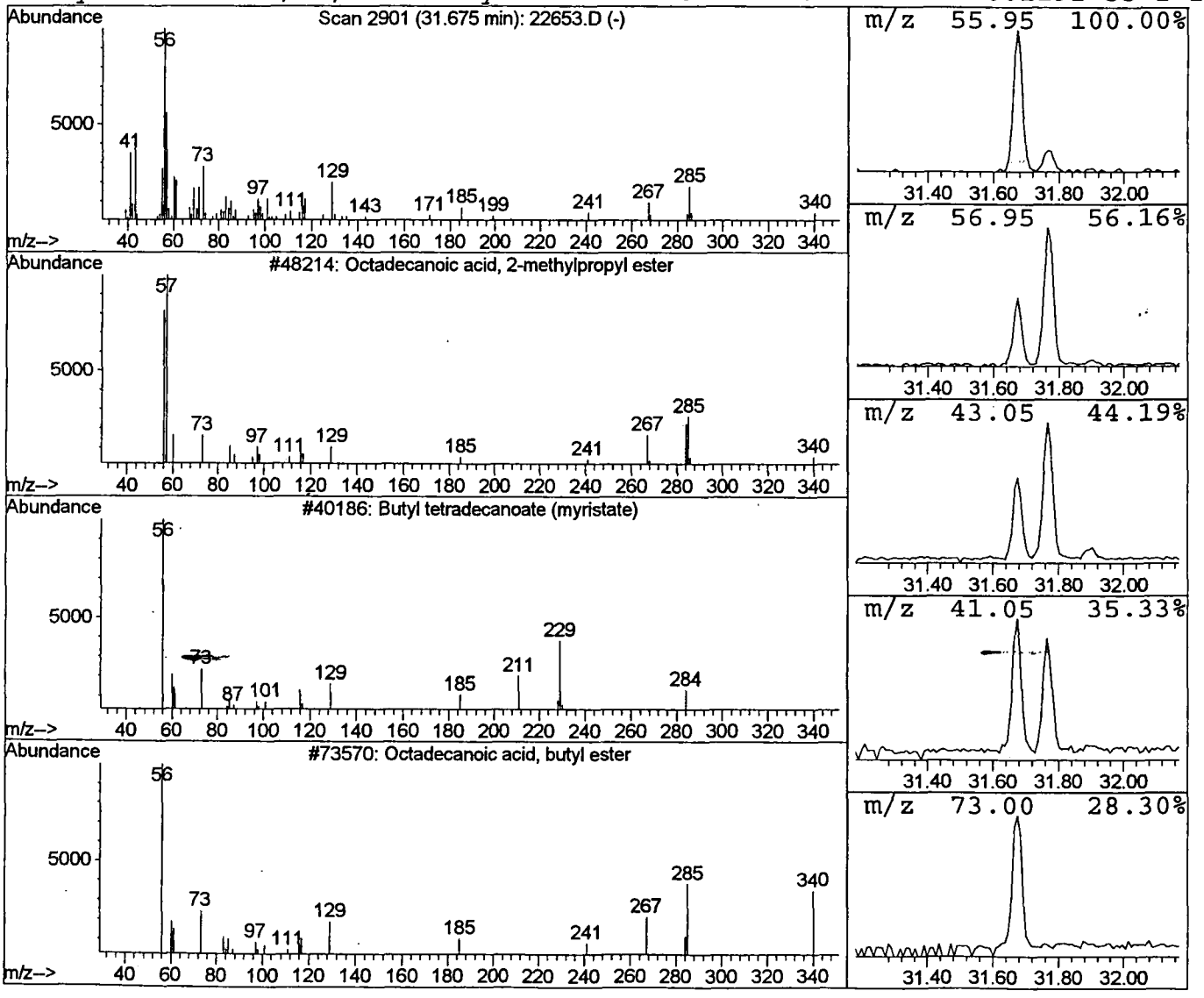
Quant Method : C:\HPCHEM\1\METHODS\NP091101.M (RTE Integrator)
Title : 209 625/TCLP calibration table 7/16/01
Library : C:\DATABASE\NBS75K.L

****

Peak Number 7 Octadecanoic acid, 2-methylpro Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
31.68	0.76 ug/l	112947	Chrysene-d12	33.16

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Octadecanoic acid, 2-methylpropyl e	340	C22H44O2	000646-13-9	83
2	Butyl tetradecanoate (myristate)	284	C18H36O2	000000-00-0	72
3	Octadecanoic acid, butyl ester	340	C22H44O2	000123-95-5	40
4	Cyclobutanone, 3,3-dimethyl-	98	C6H10O	001192-33-2	27



Library Search Compound Report

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Acq On : 28 Sep 2001 11:13 pm
Sample : gc/ms unknown
Misc :
MS Integration Params: LSCINT.P

Vial: 12
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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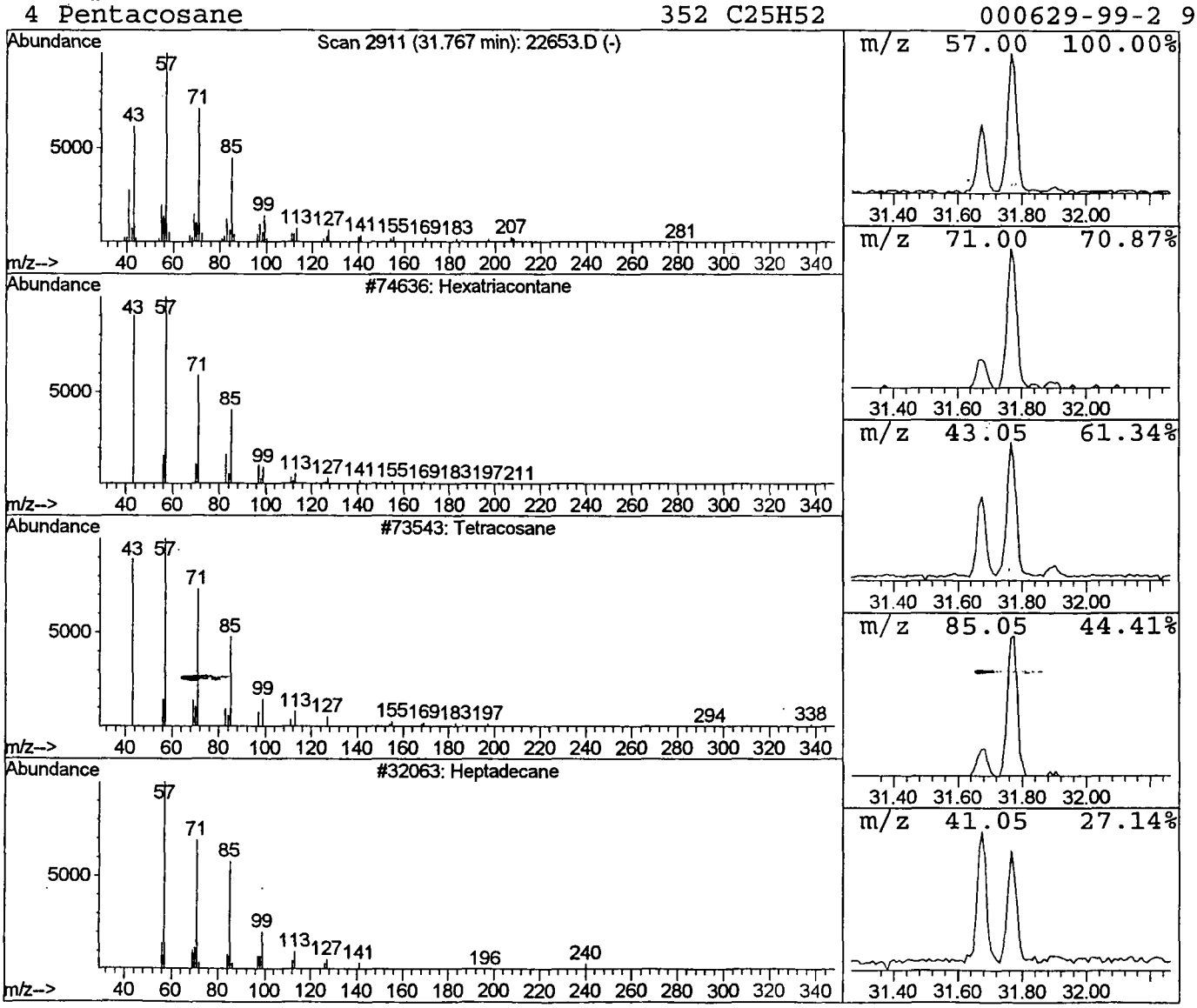
Peak Number 8 Hexatriacontane Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
31.77	0.74 ug/l	109618	Chrysene-d12	33.16

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Hexatriacontane	507	C36H74	000630-06-8	91
2	Tetracosane	338	C24H50	000646-31-1	91
3	Heptadecane	240	C17H36	000629-78-7	91
4	Pentacosane	352	C25H52	000629-99-2	90



Library Search Compound Report

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Misc :
MS Integration Params: LSCINT.P

Vial: 12
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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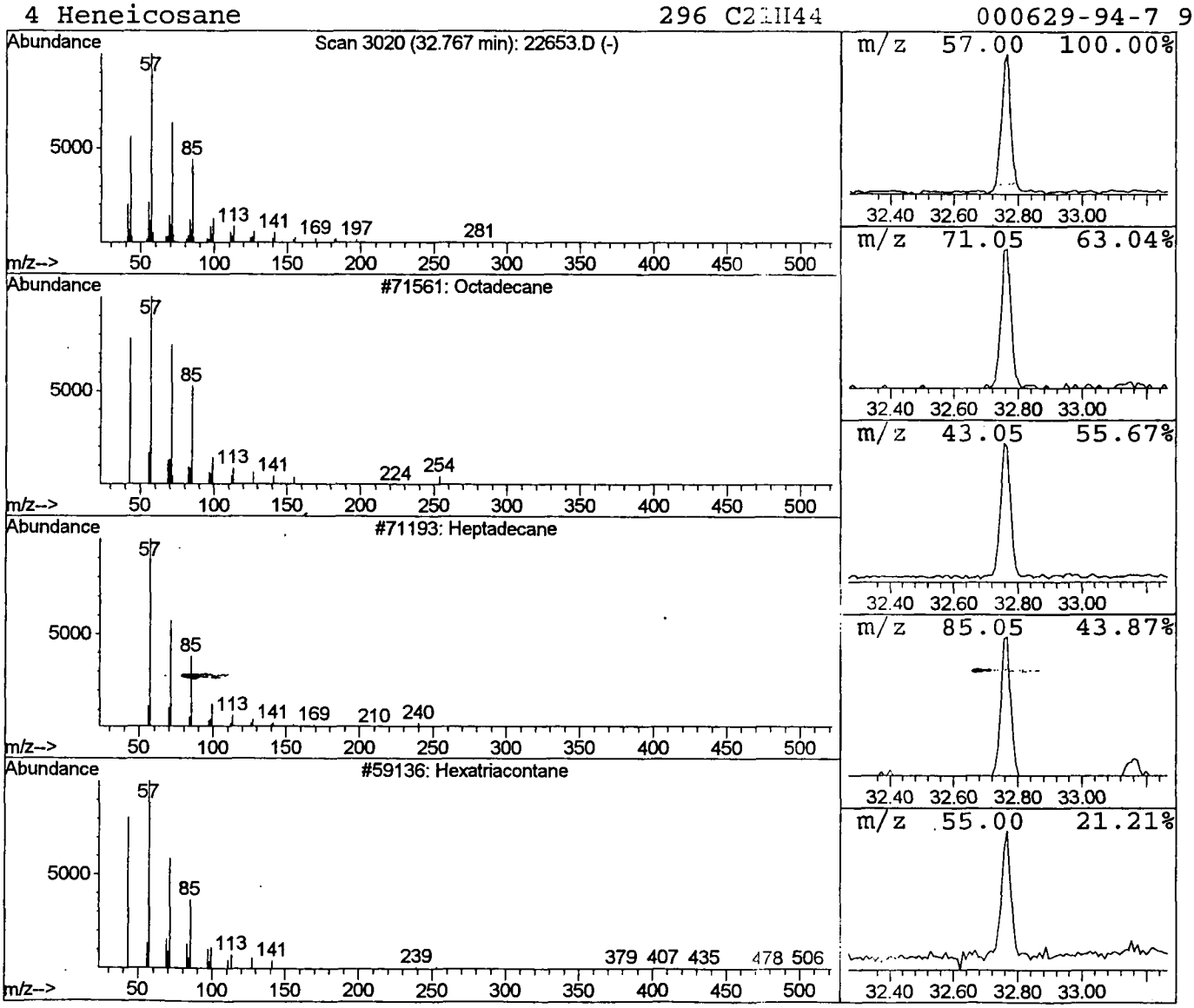
Peak Number 9 Octadecane — Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
32.77	0.78 ug/l	116242	Chrysene-d12	33.16

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Octadecane	254	C18H38	000593-45-3	91
2	Heptadecane	240	C17H36	000629-78-7	91
3	Hexatriacontane	507	C36H74	000630-06-8	91
4	Heneicosane	296	C21H44	000629-94-7	90



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Vial: 12
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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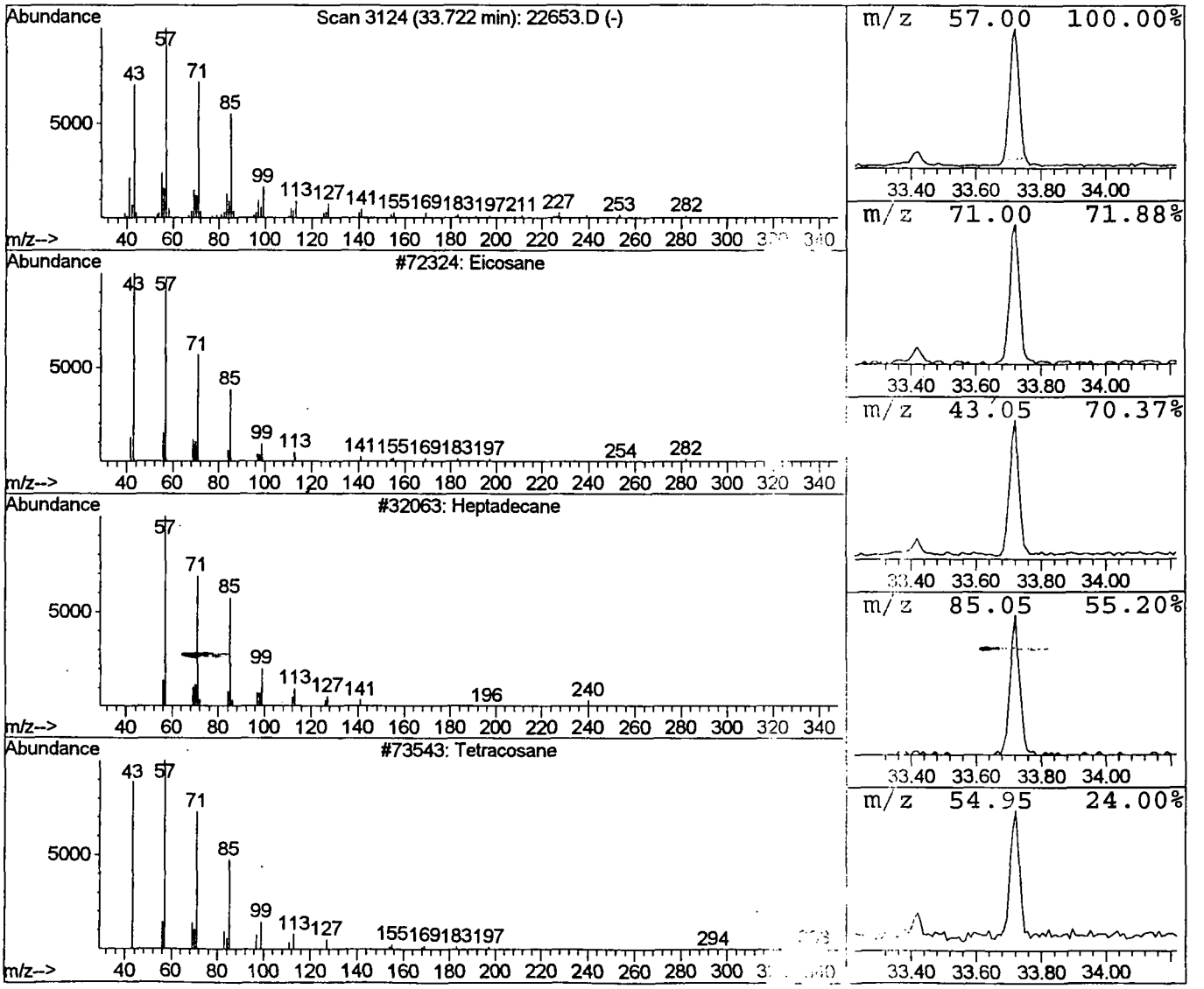
Peak Number 10 Eicosane Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
33.72	1.08 ug/l	160266	Chrysene-d12	33.16

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Eicosane	282	C20H42	000112-95-8	93
2	Heptadecane	240	C17H36	000629-78-7	91
3	Tetracosane	338	C24H50	000646-31-1	91
4	Octadecane	254	C18H38	000593-45-3	91



Library Search Compound Report

Data File : C:\HPCHEM\1\DATA\SEP2801\22653.D
Acq On : 28 Sep 2001 11:13 pm
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Misc :
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Vial: 12
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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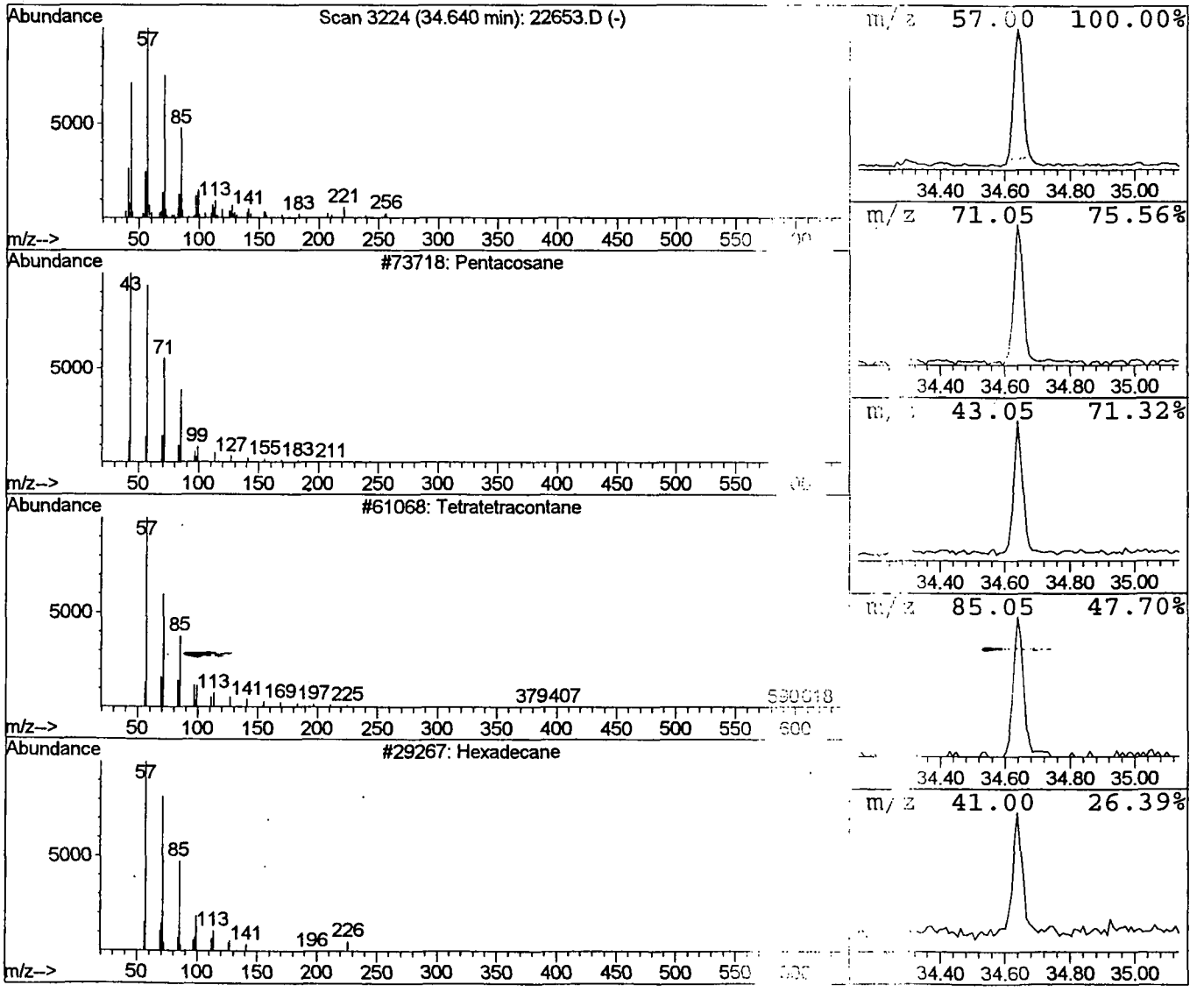
Peak Number 11 Pentacosane Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
34.64	0.75 ug/l	111676	Chrysene-d12	33.16

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Pentacosane	352	C25H52	000629-99-2	91
2	Tetratetracontane	619	C44H90	007098-22-8	91
3	Hexadecane	226	C16H34	000544-76-3	90
4	Heneicosane	296	C21H44	000629-94-7	90



Library Search Compound Report

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Vial: 12
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, 1
Library : C:\DATABASE\NBS75K.L

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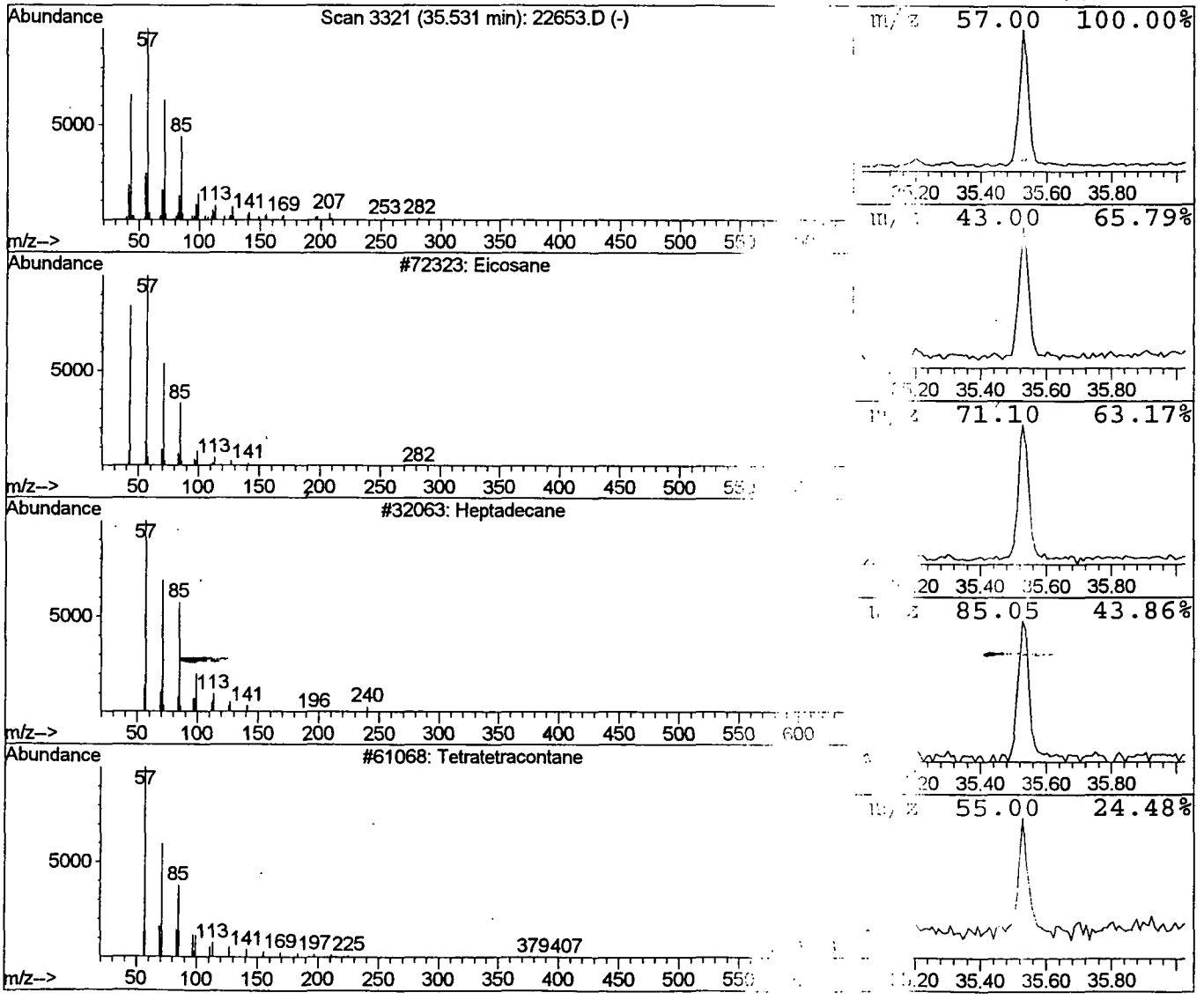
Peak Number 12 Eicosane Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
35.53	0.74 ug/l	101310	Perylene-d12	37.28

Hit# of 5	Tentative ID	MW	Molecular Formula	CAS#	Qual
1	Eicosane	282	C20H42	000112-95-8	93
2	Heptadecane	240	C17H36	000629-78-7	91
3	Tetratetracontane	619	C44H90	007098-22-8	90
4	Hexatriacontane	507	C36H74	000630-06-8	90



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Misc : Multiplr: 1.00
MS Integration Params: LSCINT.P

Quant Method : C:\HPCHEM\1\METHODS\NP091101.M (RTE Integrator)
Title : 209 625/TCLP calibration tab[illegible] 1
Library : C:\DATABASE\NBS75K.L

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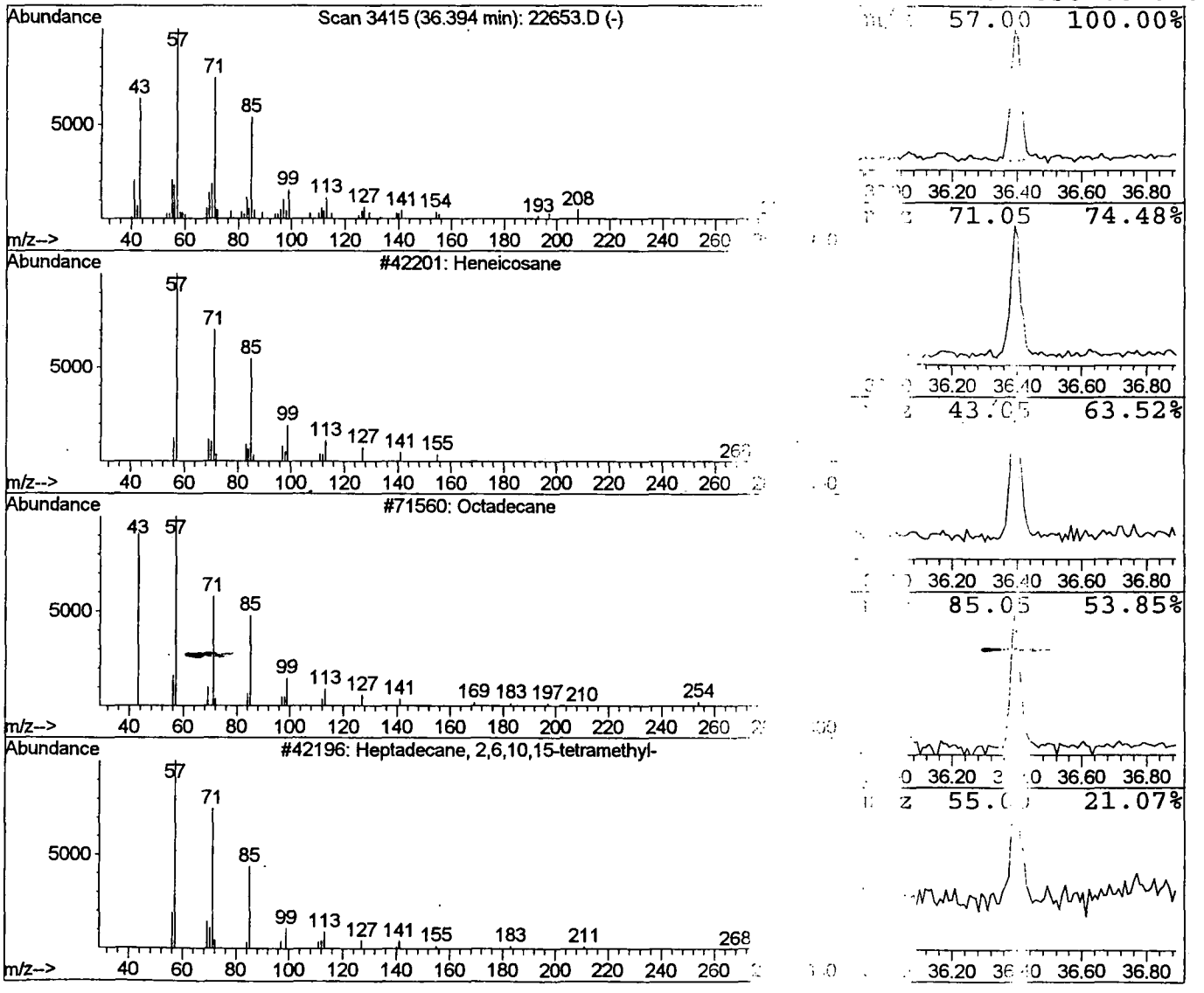
Peak Number 13 Heneicosane Concentration Rank 10

R.T.	EstConc	Area	Relati[illegible] to IS[illegible]	R.T.
36.39	0.59 ug/l	80479	Perylene-[illegible]	37.28

Hit# of 5	Tentative ID	MW	Mo[illegible]rm	CAS#	Qual
1	Heneicosane	296	C21[illegible]	000629-94-7	90
2	Octadecane	254	C[illegible]	000593-45-3	86
3	Heptadecane, 2,6,10,15-tetramethyl-	296	C2[illegible]	054833-48-6	86
4	Hexatriacontane	507	C3[illegible]	000630-06-8	86



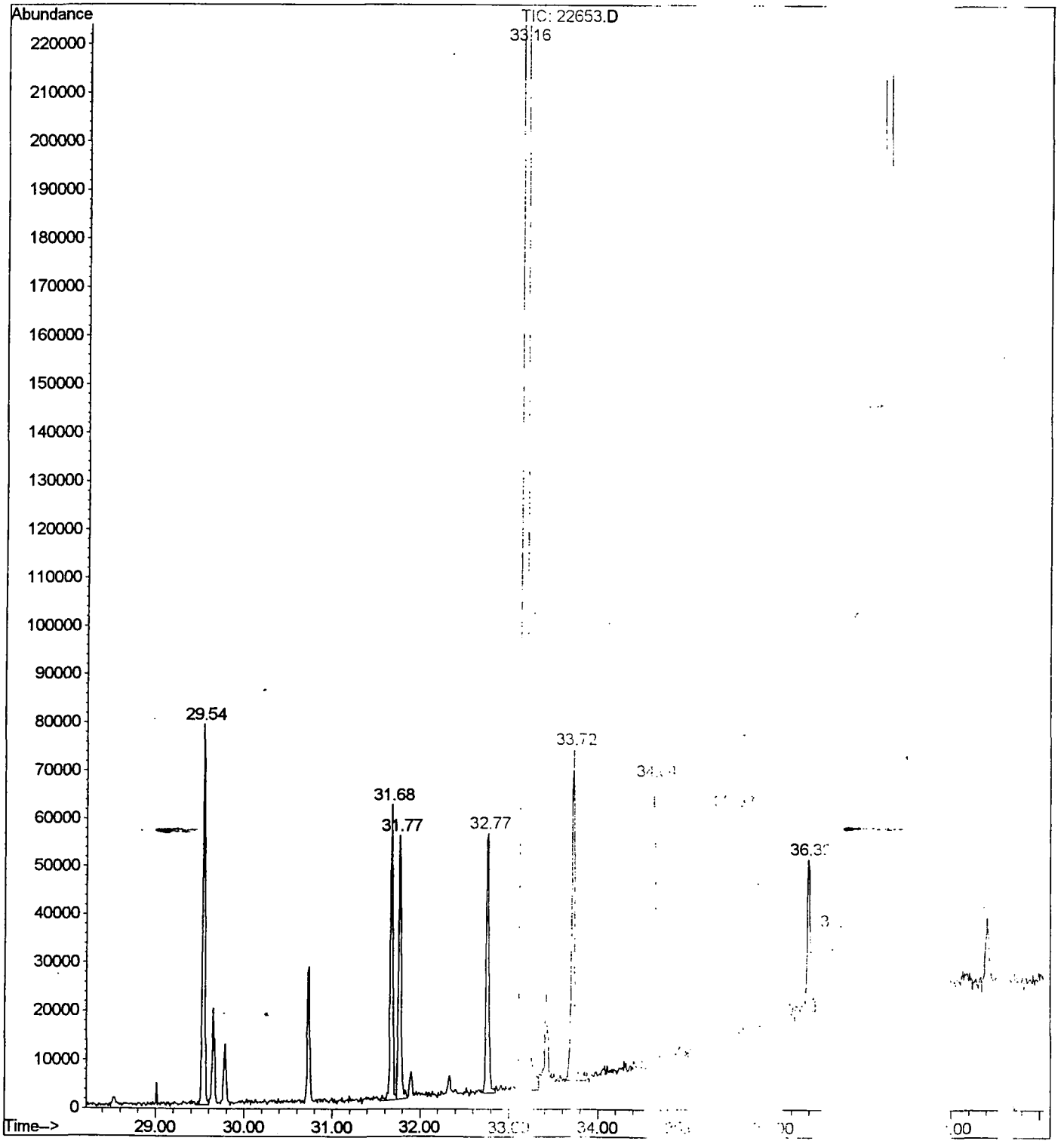
Tentatively Identified Compound (L ' ' ary

Operator ID: 209 GALOS Date Acquired: 28 Sep 1 :13 pm
Data File: C:\HPCHEM\1\DATA\SEP2801\22653.D
Name: gc/ms unknown
Misc:
Method: C:\HPCHEM\1\METHODS\NP091101.M (RTE Integ or)
Title: 209 625/TCLP calibration table 7/16/01
Library Searched: C:\DATABASE\NBS75K.L

TIC Top Hit name	RT	EstConc	Units	Area	LSt #	ISRT	ISArea	ISConc
~~3-Hexanone~~	6.39	0.6	ug/l	69191		11.79	2500960	20.0
~~2-Hexanone~~	6.49	1.1	ug/l	133042		11.79	500960	20.0
~~3-Hexanol~~	6.63	0.5	ug/l	56478		11.79	500960	20.0
2-Pentanol, 4-methyl	6.74	0.6	ug/l	74340		11.79	2500960	20.0
1,~~4-Dichlor~~obenzene-	11.64	0.5	ug/l	6711		11.79	2500960	20.0
~~Butyl hexadecanoate~~	29.54	1.0	ug/l	150248		33.16	2976970	20.0
Octadecanoic acid, 2	31.68	0.8	ug/l	112947		33.16	2976970	20.0
Hexatriacontane	31.77	0.7	ug/l	109618		33.16	2976970	20.0
Octadecane	32.77	0.8	ug/l	116242		33.16	2976970	20.0
Eicosane	33.72	1.1	ug/l	160266		33.16	2976970	20.0
Pentacosane	34.64	0.8	ug/l	111676		33.16	2976970	20.0
Eicosane	35.53	0.7	ug/l	101310		37.28	2735720	20.0
Heneicosane	36.39	0.6	ug/l	80479		37.28	2735720	20.0

22653.D NP091101.M Fri Oct 05 08:57:08 2(

File : C:\HPCHEM\1\DATA\SEP2801\22653.D
Operator : 209 GALOS
Acquired : 28 Sep 2001 11:13 pm using Acq
Instrument : GC/MS 209
Sample Name: gc/ms unknown
Misc Info :
Vial Number: 12



Comments for Sample AD22653 - Analysis $GCMS

Westchester Dept. of Labs & Research
User: Vinci, David L
Date: 10-12-2001 Time: 13:57:47

A scan of the extract prepared for the 508 analysis, from 35 - 550 amu using a mass spectrometer, does not reveal the presence of any non-target compounds generally different than those in the Laboratory Blank, except that the levels are higher in the sample than in the Laboratory Blank.