User: Galos, Marie
Westchester Dept. of Labs & Research
Date: 10/12/2001 Time: 09:10:35

Sample: AD23258
Analysis: $GCMS GCMS Scan For Unknowns
Units: ug
Started: 10/12/01 09:08 Ended: 10/12/01 09:08 Analyst: MG
Page: 1

	Component Name	Viol	Result
1	Other unknown compounds		see comment

Quantitation Report (QT Reviewed)

Data File : C:\HPCHEM\1\DATA\OCT0901\23258.D Vial: 4
Acq On : 9 Oct 2001 4:17 pm Operator: 209 GALOS
Sample : g c/ms unknown Inst : GC/MS 209
Misc : Multiplr: 1.00
MS Integration Params: RTEINT.P
Quant Time: Oct 9 17:06 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	766005	20.00	ug/l	-0.14
19) Naphthalene-d8	15.37	136	3082292	20.00	ug/l	-0.15
31) Acenaphthene-d10	20.66	164	1451293	20.00	ug/l	-0.15
49) Phenanthrene-d10	25.09	188	2063215	20.00	ug/l	-0.14
59) Chrysene-d12	33.13	240	1179723	20.00	ug/l	-0.15
68) Perylene-d12	37.25	264	793233	20.00	ug/l	-0.17
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	11.76	132	631	0.01	ug/l	0.34
Spiked Amount 75.000			Recovery =	0.01%		
7) 1,2-Dichlorobenzene-d4	12.28	152	8522	0.23	ug/l	-0.15
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	18.80	172	3259	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	29.98	244	113	0.00	ug/l	-0.17
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	11.09	94	182	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

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Misc : Multiplr: 1.00
MS Integration Params: RTEINT.P
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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
17)	N-Nitroso-di-n-propylamine	0.00	70	0	N.D.	
18)	Hexachloroethane	0.00	117	0	N.D.	
21)	Nitrobenzene	12.99	77	1107	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.	
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	20.95	165	384	N.D.	
45)	Diethylphthalate	22.16	149	399	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	25.09	178	682	N.D.	
56)	Anthracene	25.09	178	682	N.D.	
57)	Di-n-butylphthalate	27.09	149	7155	N.D.	
58)	Fluoranthene	0.00	202	0	N.D.	
60)	3,3'-Dichlorobenzidine	0.00	252	0	N.D.	

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Quantitation Report (QT Reviewed)

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Acq On : 9 Oct 2001 4:17 pm Operator: 209 GALOS
Sample : g c/ms unknown Inst : GC/MS 209
Misc : Multiplr: 1.00
MS Integration Params: RTEINT.P
Quant Time: Oct 9 17:06 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
61)	Benzidine	0.00	184	0	N.D.	
63)	Pyrene	0.00	202	0	N.D.	
64)	Butylbenzylphthalate	31.59	149	1376	N.D.	
65)	Benzo(a)anthracene	33.14	228	2973	N.D.	
66)	Bis(2-ethylhexyl)phthalate	33.38	149	10072	N.D.	
67)	Chrysene	33.20	228	177	N.D.	
69)	Di-n-Octylphthalate	35.15	149	2029	N.D.	
70)	Benzo(b)fluoranthene	36.14	252	280	N.D.	
71)	Benzo(k)fluoranthene	36.14	252	280	N.D.	
72)	Benzo(a)pyrene	37.25	252	3085	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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Acq On : 9 Oct 2001 4:17 pm Operator: 209 GALOS
Sample : g c/ms unknown Inst : GC/MS 209
Misc : Multiplr: 1.00
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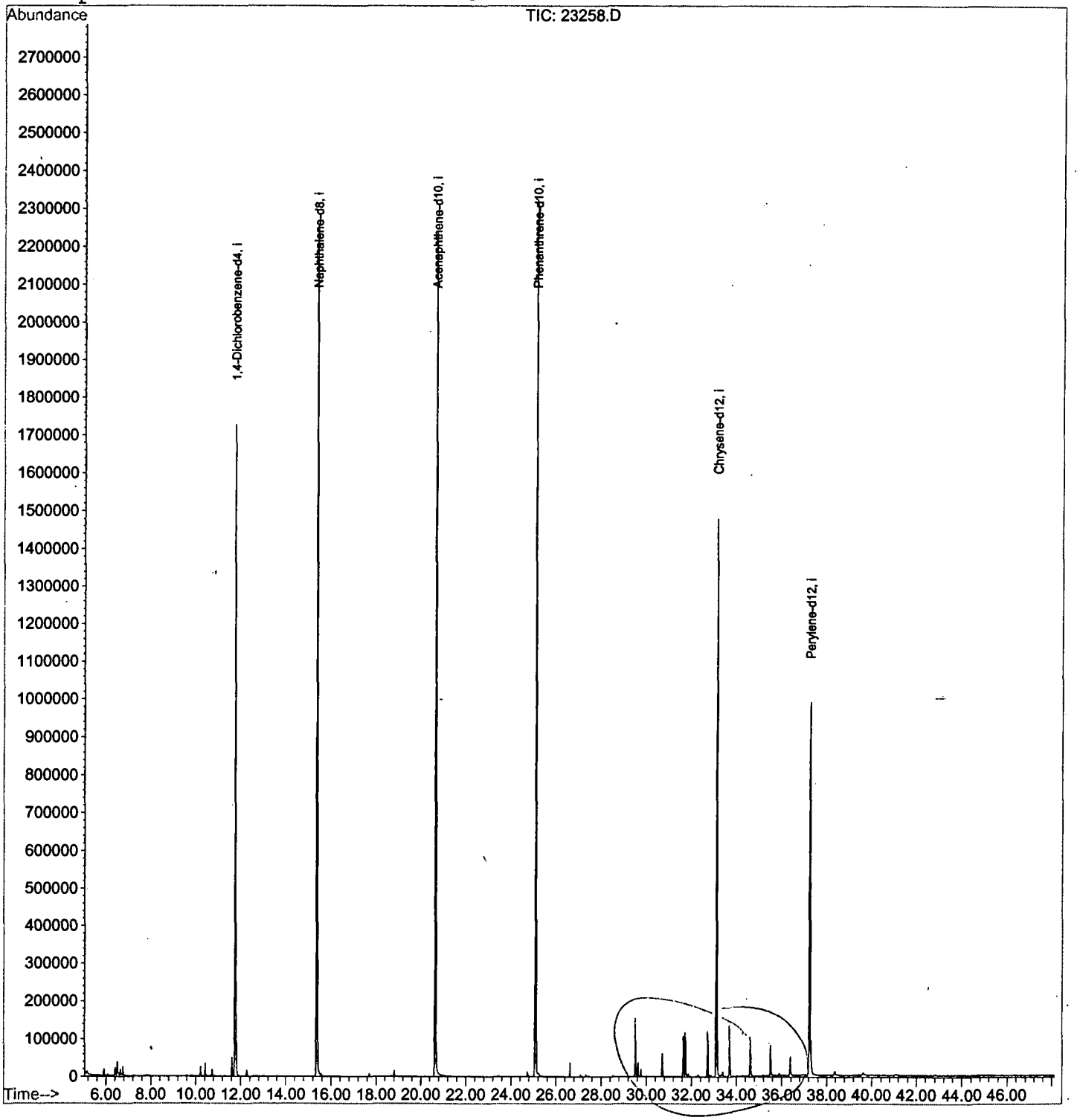
Title : 209 625/TCLP calibration table 7/16/01
Last Update : Thu Sep 13 12:47:04 2001
Response via : Initial Calibration



LSC Area Percent Report

Data File : C:\HPCHEM\1\DATA\OCT0901\232[illegible].D Vial: 4
Acq On : 9 Oct 2001 4:17 pm [illegible]erator: 209 GALO
Sample : g c/ms unknown [illegible]t : GC/MS 209
Misc : [illegible]tiplr: 1.00
MS Integration Params: LSCINT.P

Method : C:\HPCHEM\1\METHODS\NP091101.M (RTE Int [illegible]tor)
Title : 209 625/TCLP calibration table 7/16/01
Smoothing : OFF Filterin
Sampling : 1 Min Are % of largest Peak
Start Thrs: 0.2 Max Peak [illegible]00
Stop Thrs : 0 Peak Locatio [illegible]DP

If leading or trailing edge < 100 prefer < Baseline > else tangent >
Peak separation: 5

Signal : TIC

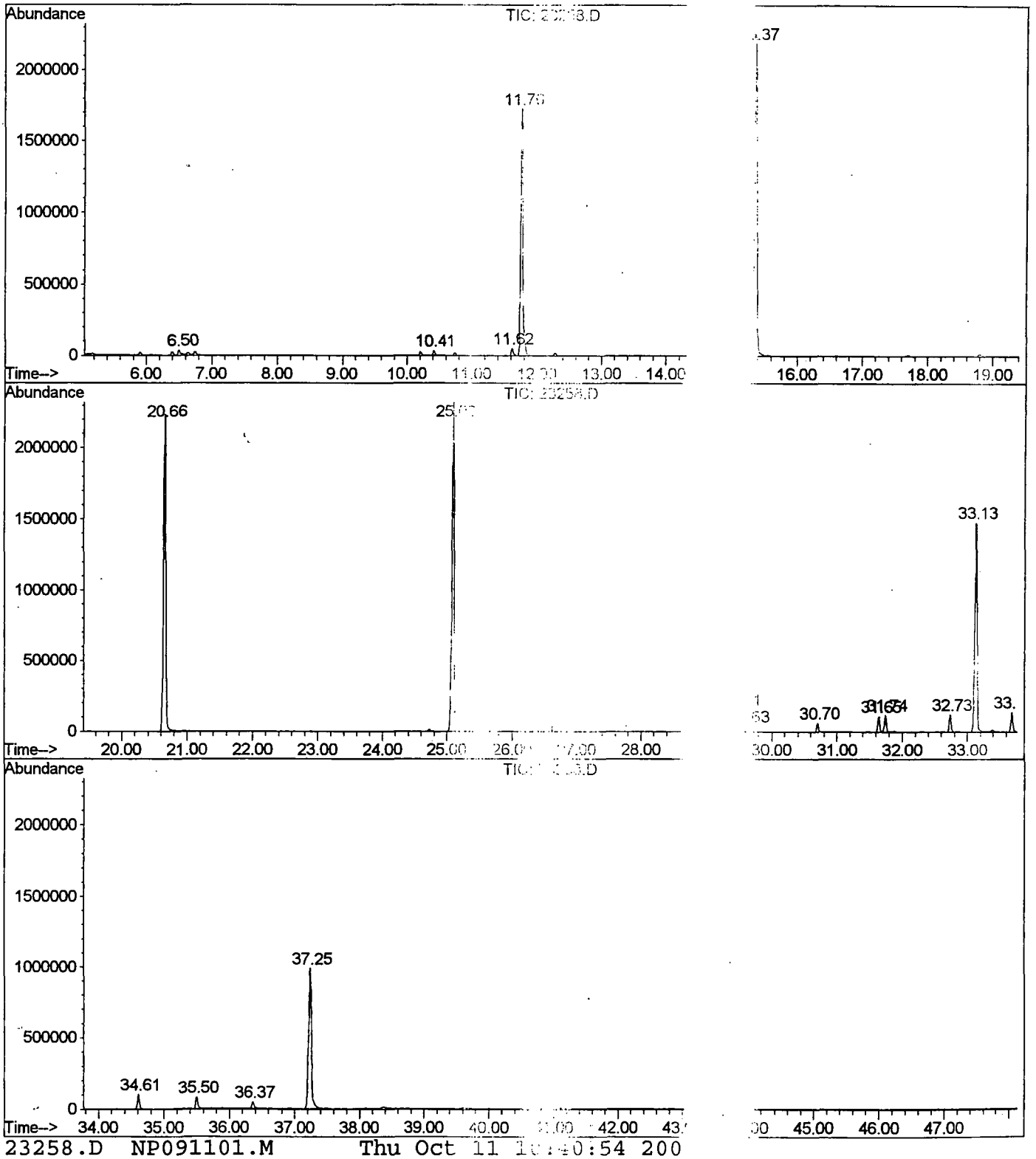
peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	peak area	%	% of total
1	6.504	155	159	168	rVV	36800	85219	[illegible]	0.286%
2	10.415	581	585	594	rBV	34239	68420	[illegible]	0.230%
3	11.617	711	716	726	rBV	49961	110634	[illegible]	0.372%
4	11.764	726	732	748	rBV	1726590	[illegible]005076	6	13.460%
5	15.372	1117	1125	1142	rBV	2256[illegible]56	[illegible]653363	9	19.000%
6	20.660	1692	1701	1721	rBV	2235220	5809313	10	19.524%
7	25.094	2173	2184	2195	rBV2	2322371	5295156	9	17.796%
8	29.510	2660	2665	2672	rBV2	153786	295775	[illegible]	0.994%
9	29.629	2672	2678	2685	rVB	37407	76204	[illegible]	0.256%
10	30.703	2788	2795	2800	rBV	61470	122422	[illegible]	0.411%
11	31.649	2893	2898	2903	rBV	105336	210510	[illegible]	0.707%
12	31.741	2903	2908	2913	rBV	115646	[illegible]3828	[illegible]	0.752%
13	32.732	3011	3016	3023	rBV	118[illegible]67	[illegible]40258	[illegible]	0.807%
14	33.127	3051	3059	3073	rBV2	1480[illegible]09	[illegible]707542	6	12.460%
15	33.687	3113	3120	3133	rBV	135[illegible]82	[illegible]6357	[illegible]	0.962%
16	34.614	3216	3221	3231	rVB	101242	213473	[illegible]	0.717%
17	35.504	3312	3318	3328	rBV	79962	183151	[illegible]	0.616%
18	36.367	3405	3412	3424	rBV2	48722	123851	[illegible]	0.416%
19	37.249	3495	3508	3523	rBV	987[illegible]87	3044414	5	10.232%

Sum of corrected areas: 29[illegible]56

23258.D NP091101.M Thu Oct 11 [illegible]0:40:[illegible] 2001

LSC Report - Integrated Chromat

File : C:\HPCHEM\1\DATA\OCT0901\23258.D
Operator : 209 GALOS
Acquired : 9 Oct 2001 4:17 pm using AcqMeth 091101
Instrument : GC/MS 209
Sample Name: g c/ms unknown
Misc Info :
Vial Number: 4
Quant File :NP091101.RES (RTE Integrator)



23258.D NP091101.M Thu Oct 11 10:40:54 200

Library Search Compound Report

Data File : C:\HPCHEM\1\DATA\OCT0901\23258.D
Acq On : 9 Oct 2001 4:17 pm
Sample : g c/ms unknown
Misc :
MS Integration Params: LSCINT.P

Vial: 4
Operator: 209 GALOS
Inst : GC/MS 209
Multiplr: 1.00

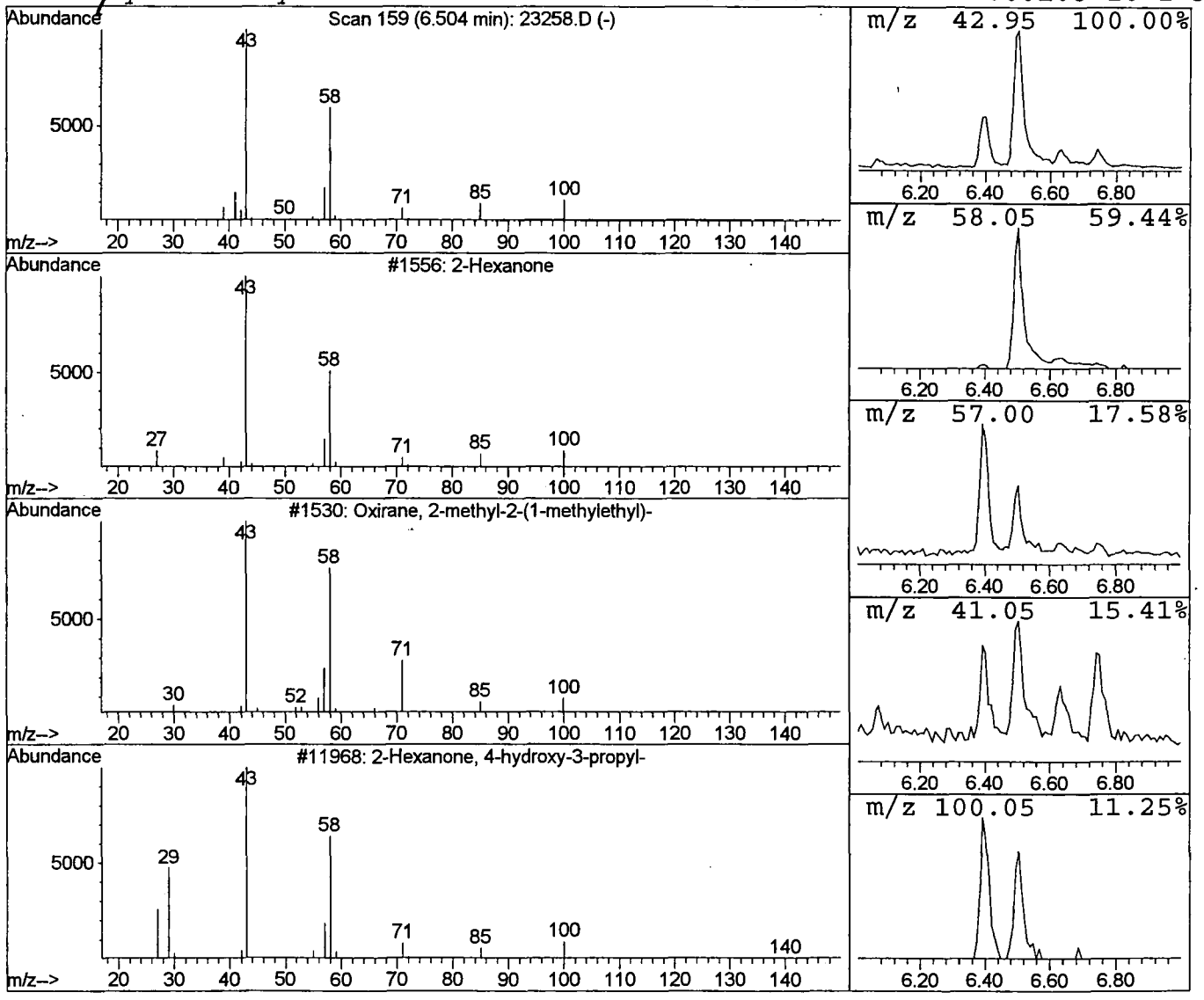
Quant Method : C:\HPCHEM\1\METHODS\NP091101.M (RTE Integrator)
Title : 209 625/TCLP calibration table 7/16/01
Library : C:\DATABASE\NBS75K.L

****

Peak Number 1 2-Hexanone Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.50	0.43 ug/l	85219	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	78
3	2-Hexanone, 4-hydroxy-3-propyl-	158	C9H18O2	062338-17-4	74
4	Methyl Isobutyl Ketone	100	C6H12O	000108-10-1	58



Library Search Compound Report

Data File : C:\HPCHEM\1\DATA\OCT0901\23258.D
Acq On : 9 Oct 2001 4:17 pm
Sample : g c/ms unknown
Misc :
MS Integration Params: LSCINT.P

Vial: 4
Operator: 209 GALOS
Inst : GC/MS 209
Multiplr: 1.00

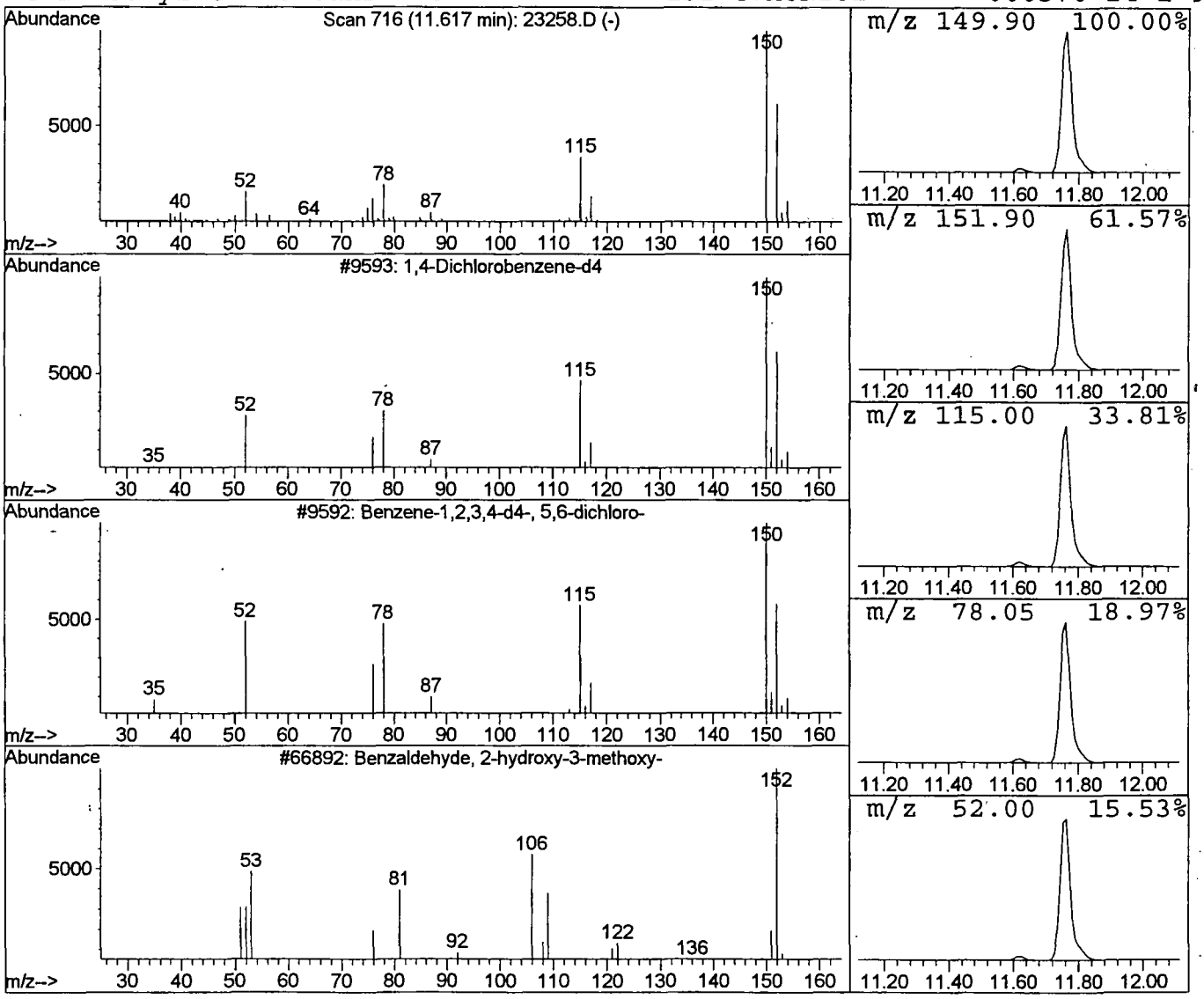
Quant Method : C:\HPCHEM\1\METHODS\NP091101.M (RTE Integrator)
Title : 209 625/TCLP calibration table 7/16/01
Library : C:\DATABASE\NBS75K.L

****

Peak Number 2 1,4-Dichlorobenzene-d4 Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.62	0.55 ug/l	110634	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	53
2	Benzene-1,2,3,4-d4-, 5,6-dichloro-	150	C6Cl2D4	002199-69-1	42
3	Benzaldehyde, 2-hydroxy-3-methoxy-	152	C8H8O3	000148-53-8	10
4	2-Methyl-6-nitroaniline	152	C7H8N2O2	000570-24-1	9



Library Search Compound Report

Data File : C:\HPCHEM\1\DATA\OCT0901\23258.D
Acq On : 9 Oct 2001 4:17 pm
Sample : g c/ms unknown
Misc :
MS Integration Params: LSCINT.P

Vial: 4
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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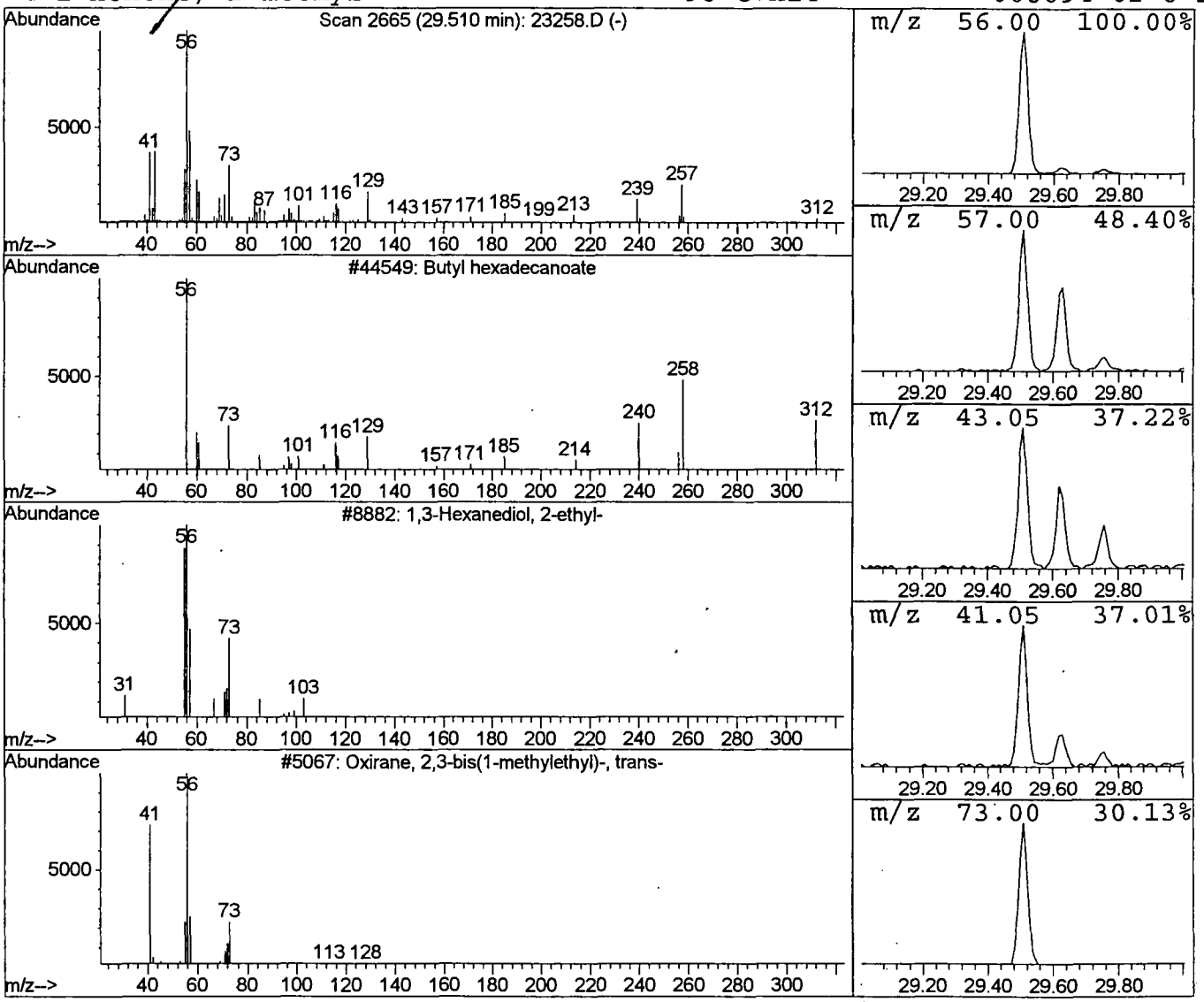
Peak Number 3 Butyl hexadecanoate Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
29.51	1.60 ug/l	295775	Chrysene-d12	33.13

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Butyl hexadecanoate	312	C20H40O2	000000-00-0	38
2	1,3-Hexanediol, 2-ethyl-	146	C8H18O2	000094-96-2	32
3	Oxirane, 2,3-bis(1-methylethyl)-, t	128	C8H16O	054644-32-5	32
4	1-Hexene, 2-methyl-	98	C7H14	006094-02-6	27



Library Search Compound Report

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Acq On : 9 Oct 2001 4:17 pm
Sample : g c/ms unknown
Misc :
MS Integration Params: LSCINT.P

Vial: 4
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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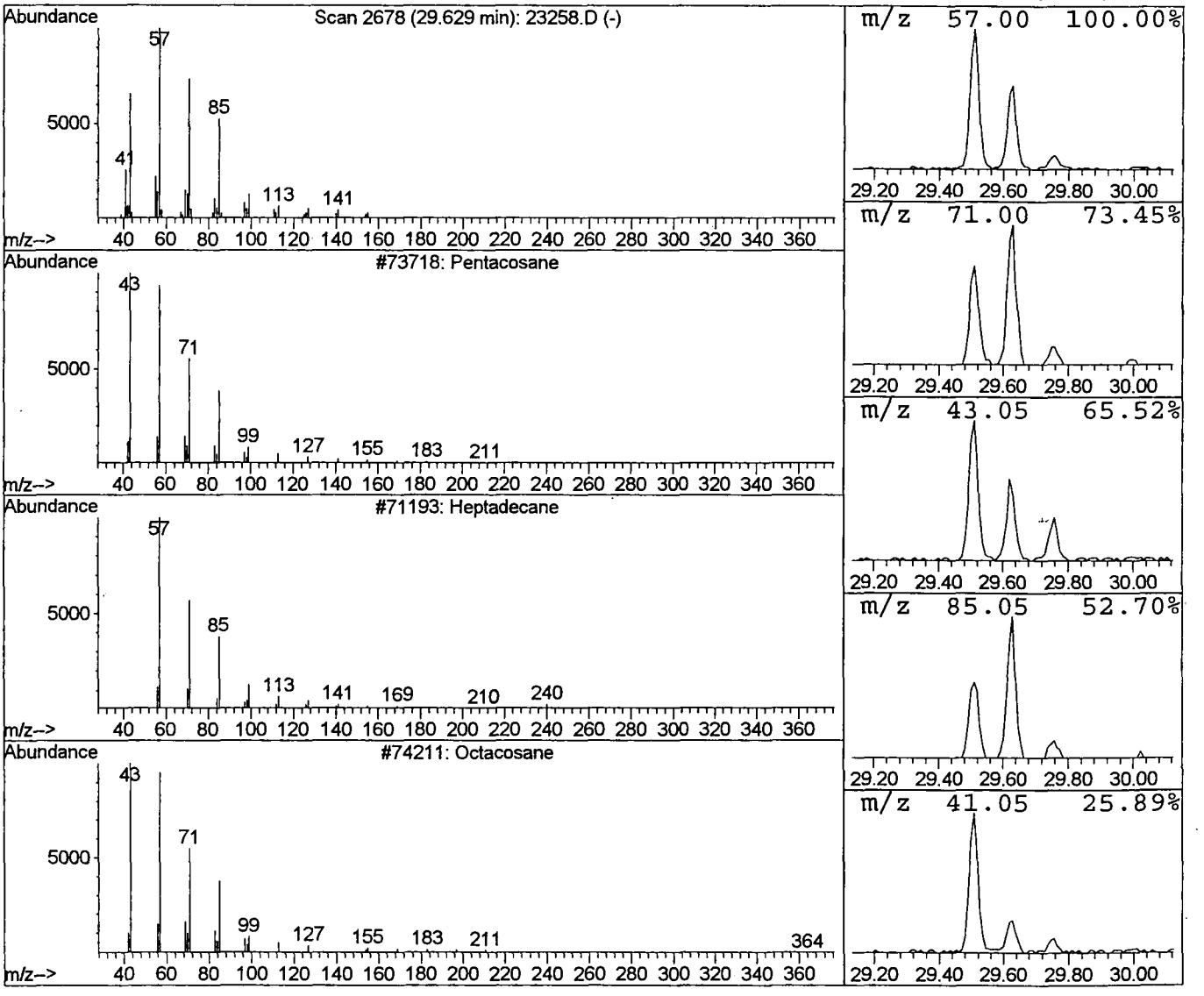
Peak Number 4 Pentacosane — Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
29.63	0.41 ug/l	76204	Chrysene-d12	33.13

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Pentacosane	352	C25H52	000629-99-2	91
2	Heptadecane	240	C17H36	000629-78-7	91
3	Octacosane	394	C28H58	000630-02-4	90
4	Docosane	310	C22H46	000629-97-0	90



Library Search Compound Report

Data File : C:\HPCHEM\1\DATA\OCT0901\23258.D
Acq On : 9 Oct 2001 4:17 pm
Sample : g c/ms unknown
Misc :
MS Integration Params: LSCINT.P

Vial: 4
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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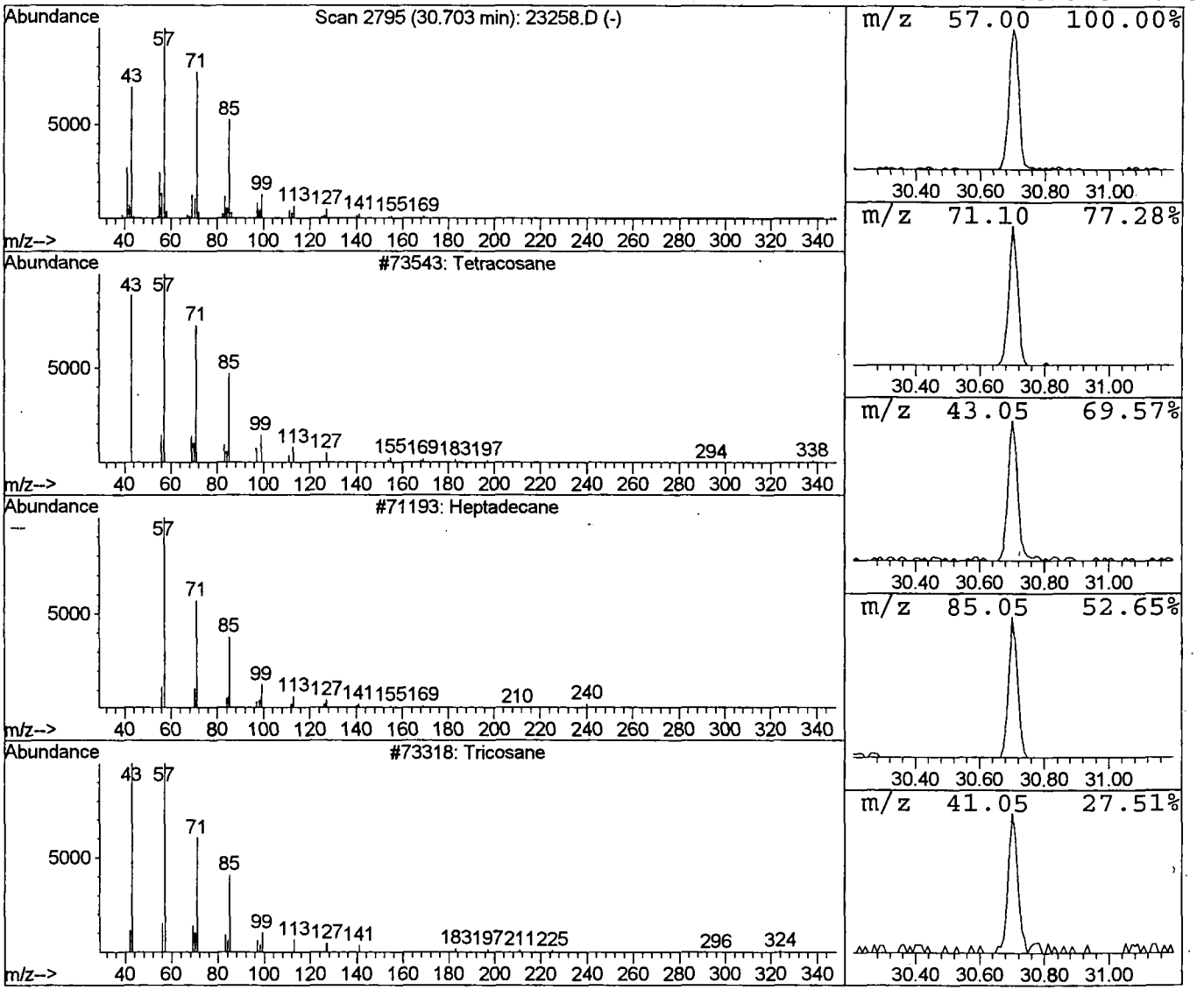
Peak Number 5 Tetracosane Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
30.70	0.66 ug/l	122422	Chrysene-d12	33.13

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Tetracosane	338	C24H50	000646-31-1	91
2	Heptadecane	240	C17H36	000629-78-7	91
3	Tricosane	324	C23H48	000638-67-5	90
4	Docosane	310	C22H46	000629-97-0	86



Library Search Compound Report

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Acq On : 9 Oct 2001 4:17 pm
Sample : g c/ms unknown
Misc :
MS Integration Params: LSCINT.P

Vial: 4
Operator: 209 GALOS
Inst : GC/MS 209
Multiplr: 1.00

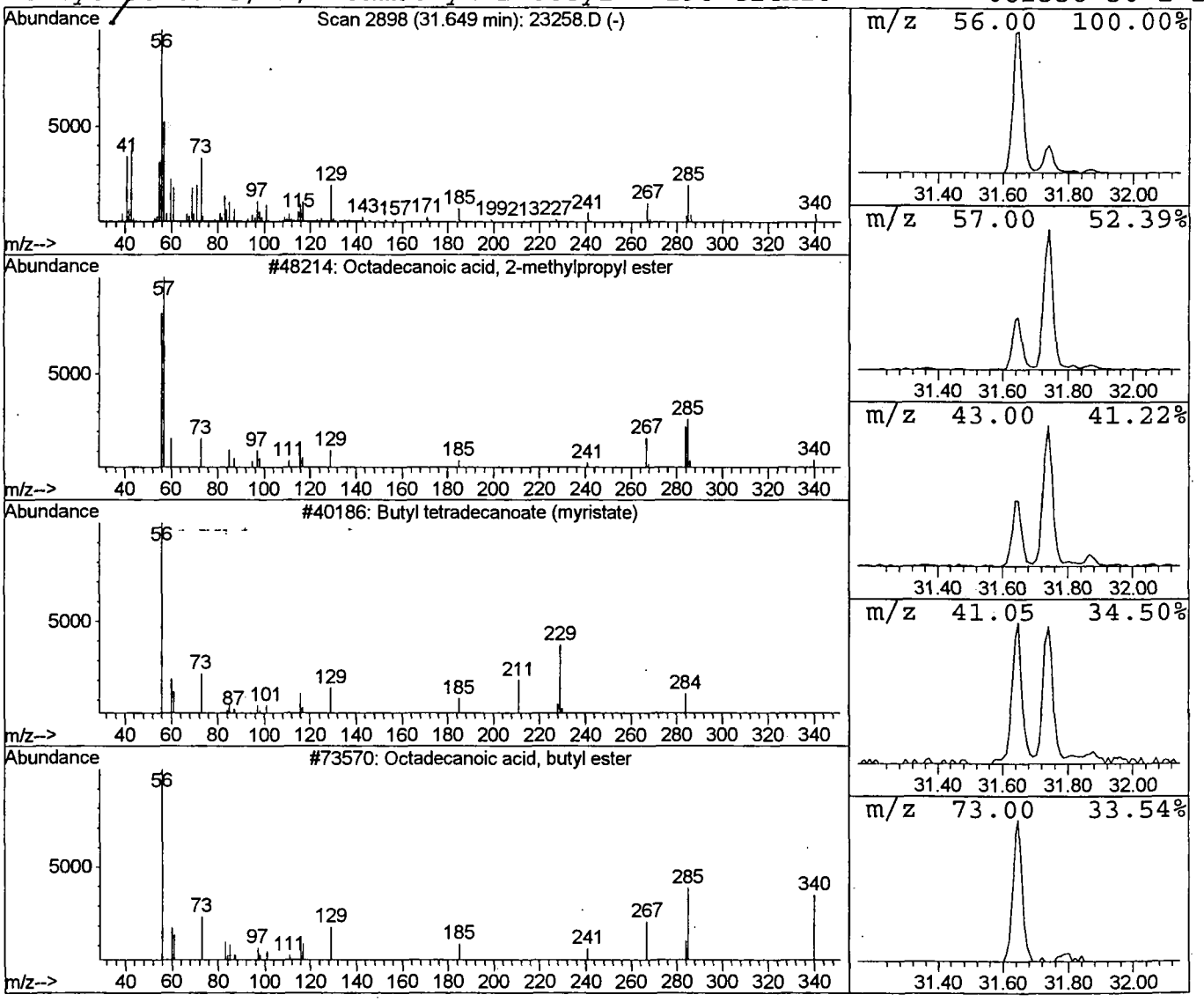
Quant Method : C:\HPCHEM\1\METHODS\NP091101.M (RTE Integrator)
Title : 209 625/TCLP calibration table 7/16/01
Library : C:\DATABASE\NBS75K.L

****

Peak Number 6 Octadecanoic acid, 2-methylpro Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
31.65	1.14 ug/l	210510	Chrysene-d12	33.13

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Octadecanoic acid, 2-methylpropyl e	340	C22H44O2	000646-13-9	64
2	Butyl tetradecanoate (myristate)	284	C18H36O2	000000-00-0	64
3	Octadecanoic acid, butyl ester	340	C22H44O2	000123-95-5	38
4	Cyclobutane, 1,1-dimethyl-2-octyl-	196	C14H28	062338-30-1	27



Library Search Compound Report

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Acq On : 9 Oct 2001 4:17 pm
Sample : g c/ms unknown
Misc :
MS Integration Params: LSCINT.P

Vial: 4
Operator: 209 GALOS
Inst : GC/MS 209
Multiplr: 1.00

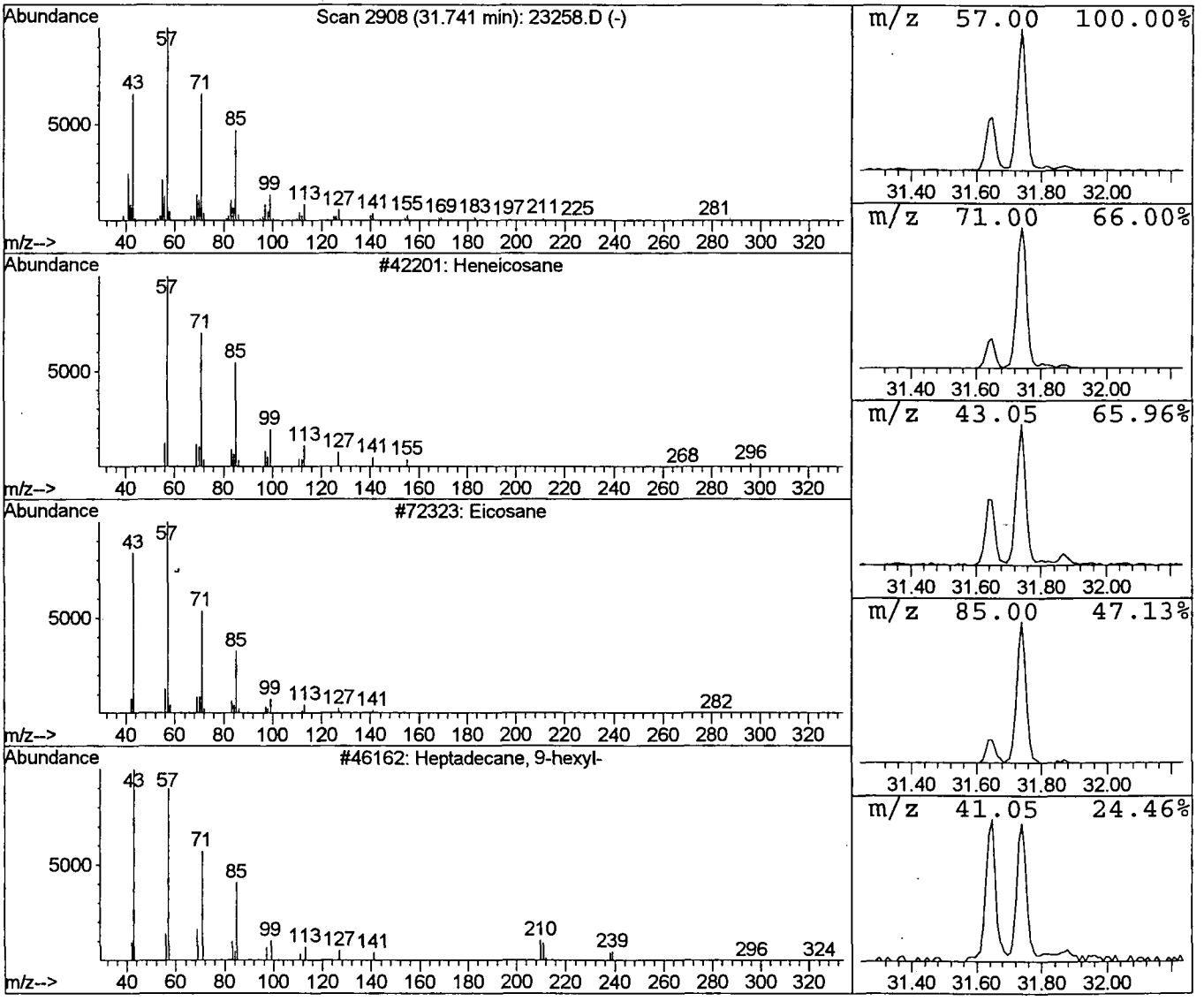
Quant Method : C:\HPCHEM\1\METHODS\NP091101.M (RTE Integrator)
Title : 209 625/TCLP calibration table 7/16/01
Library : C:\DATABASE\NBS75K.L

****

Peak Number 7 Heneicosane — Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
31.74	1.21 ug/l	223828	Chrysene-d12	33.13

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heneicosane	296	C21H44	000629-94-7	91
2	Eicosane	282	C20H42	000112-95-8	91
3	Heptadecane, 9-hexyl-	324	C23H48	055124-79-3	91
4	Pentacosane	352	C25H52	000629-99-2	91



Library Search Compound Report

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Acq On : 9 Oct 2001 4:17 pm
Sample : g c/ms unknown
Misc :
MS Integration Params: LSCINT.P

Vial: 4
Operator: 209 GALOS
Inst : GC/MS 209
Multiplr: 1.00

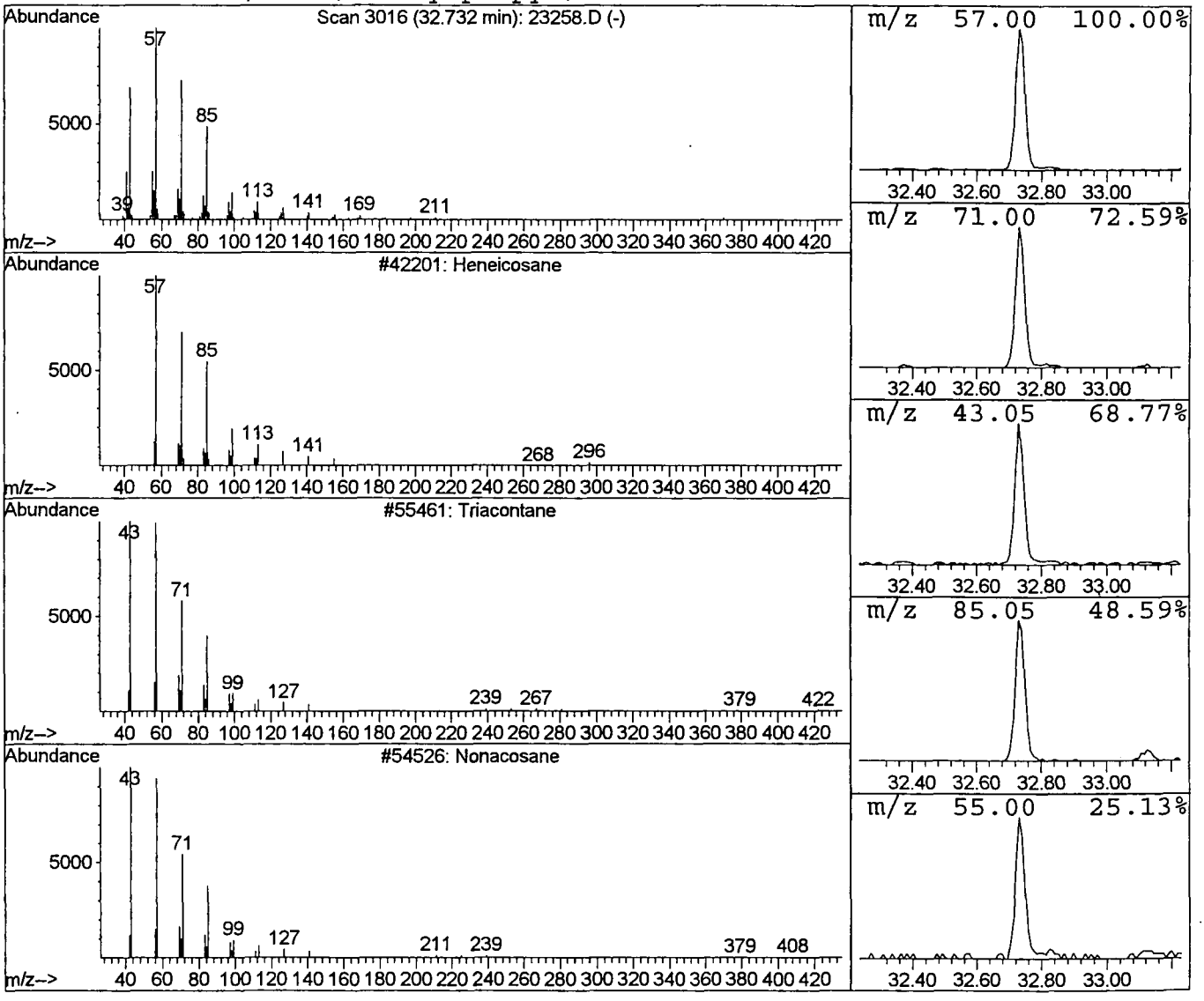
Quant Method : C:\HPCHEM\1\METHODS\NP091101.M (RTE Integrator)
Title : 209, 625/TCLP calibration table 7/16/01
Library : C:\DATABASE\NBS75K.L

****

Peak Number 8 Heneicosane Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
32.73	1.30 ug/l	240258	Chrysene-d12	33.13

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heneicosane	296	C21H44	000629-94-7	91
2	Triacontane	422	C30H62	000638-68-6	91
3	Nonacosane	408	C29H60	000630-03-5	91
4	Heneicosane, 11-(1-ethylpropyl)-	366	C26H54	055282-11-6	91



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Misc :
MS Integration Params: LSCINT.P

Vial: 4
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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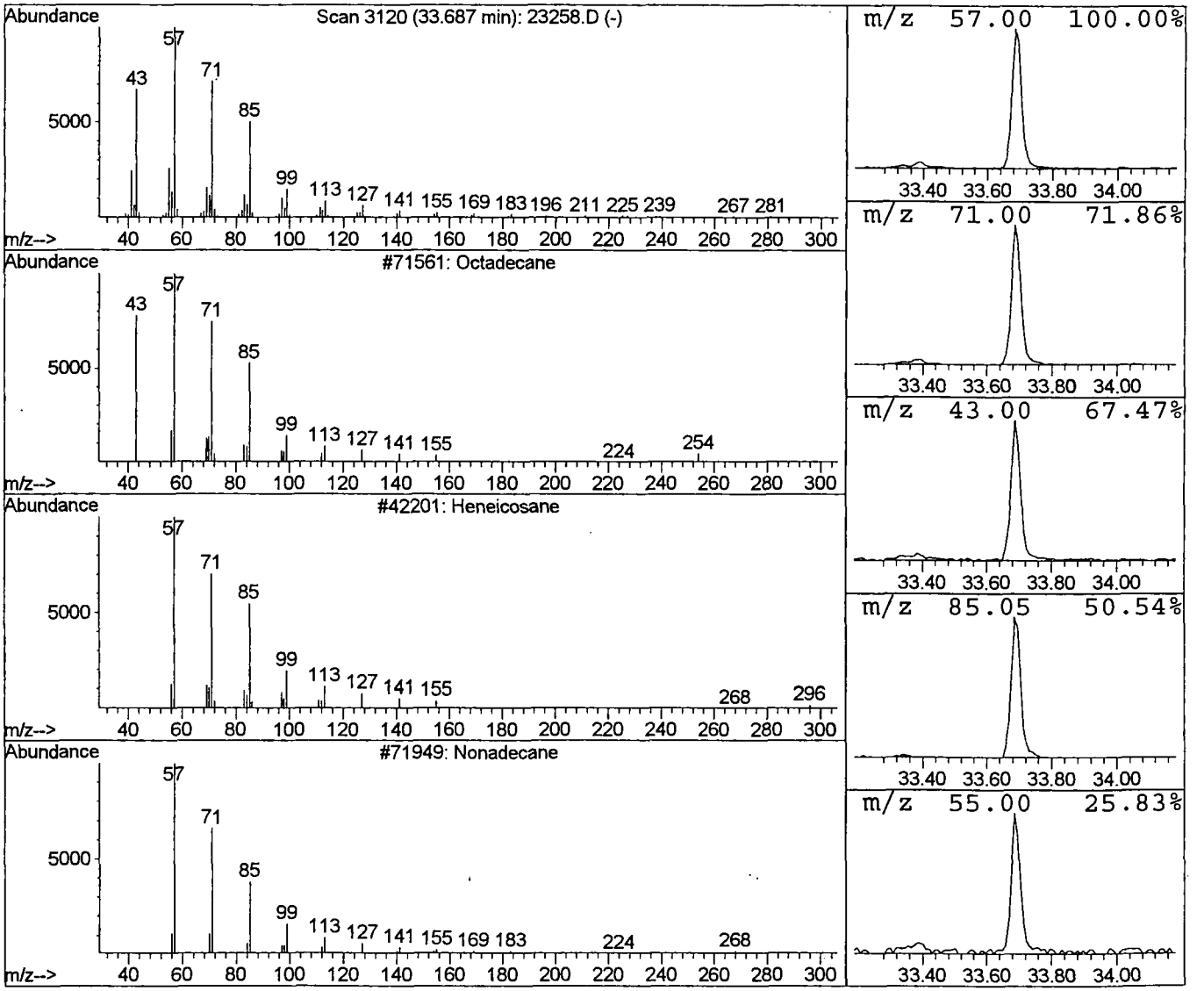
Peak Number 9 Octadecane — Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
33.69	1.54 ug/l	286357	Chrysene-d12	33.13

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Octadecane	254	C18H38	000593-45-3	91
2	Heneicosane	296	C21H44	000629-94-7	91
3	Nonadecane	268	C19H40	000629-92-5	91
4	Tricosane	324	C23H48	000638-67-5	91



Library Search Compound Report

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Acq On : 9 Oct 2001 4:17 pm
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Misc :
MS Integration Params: LSCINT.P

Vial: 4
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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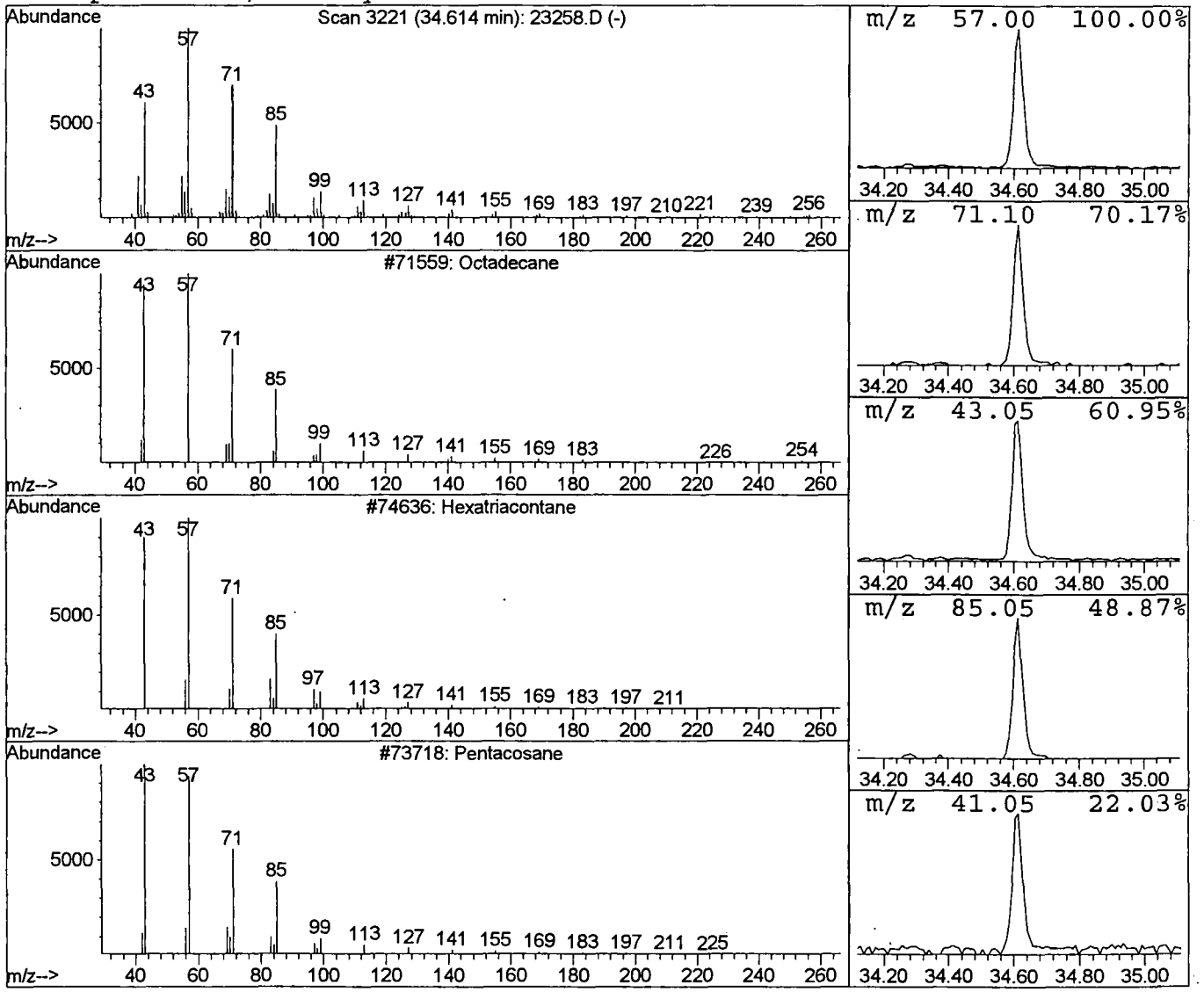
Peak Number 10 Octadecane — Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
34.61	1.15 ug/l	213473	Chrysene-d12	33.13

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Octadecane	254	C18H38	000593-45-3	94
2	Hexatriacontane	507	C36H74	000630-06-8	94
3	Pentacosane	352	C25H52	000629-99-2	91
4	Heptadecane, 9-octyl-	352	C25H52	007225-64-1	91



Library Search Compound Report

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Sample : g c/ms unknown
Misc :
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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

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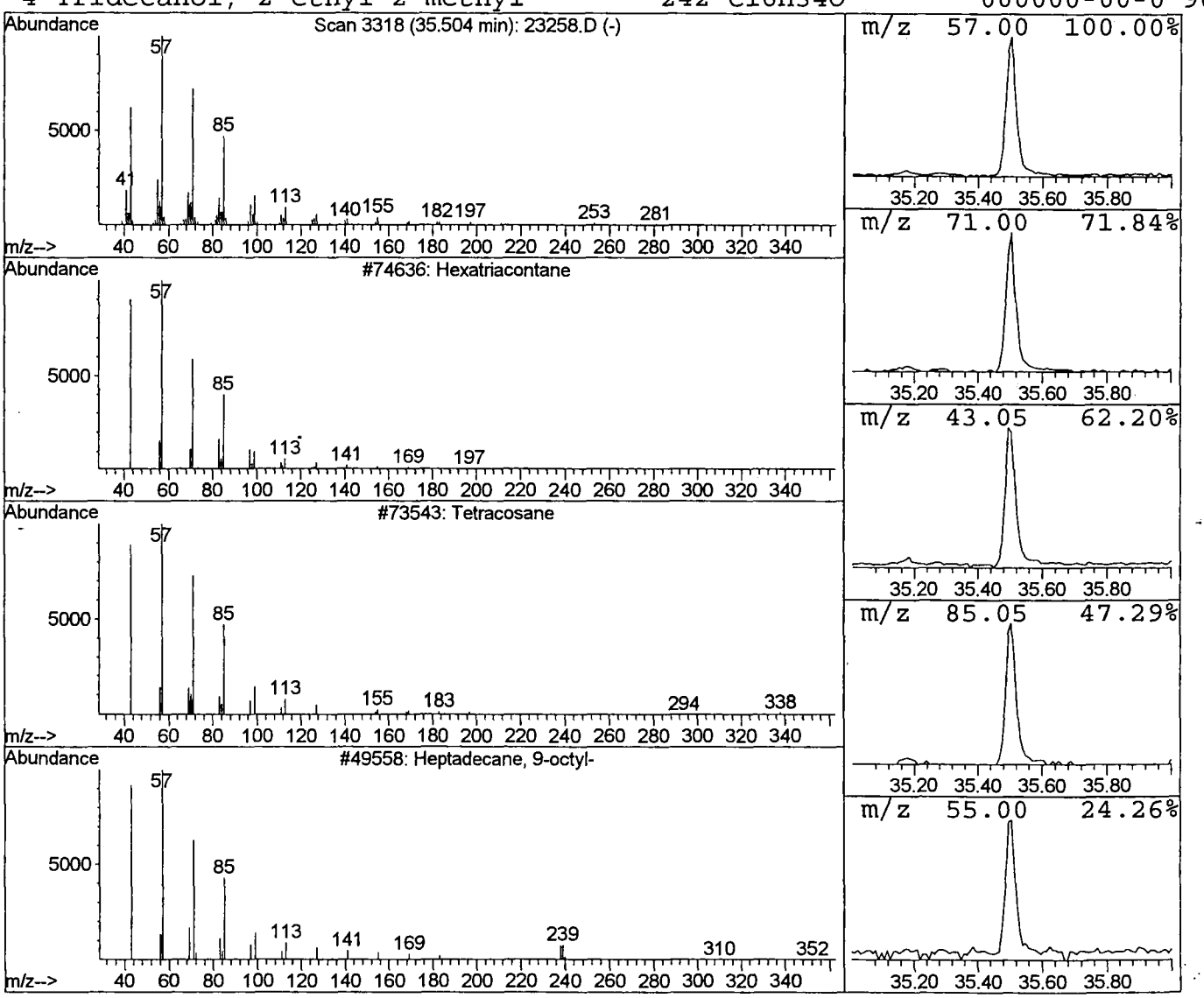
Peak Number 11 Hexatriacontane Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
35.50	1.20 ug/l	183151	Perylene-d12	37.25

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Hexatriacontane	507	C36H74	000630-06-8	94
2	Tetracosane	338	C24H50	000646-31-1	91
3	Heptadecane, 9-octyl-	352	C25H52	007225-64-1	91
4	Tridecanol, 2-ethyl-2-methyl-	242	C16H34O	000000-00-0	90



Library Search Compound Report

Data File : C:\HPCHEM\1\DATA\OCT0901\23258.D
Acq On : 9 Oct 2001 4:17 pm
Sample : g c/ms unknown
Misc :
MS Integration Params: LSCINT.P

Vial: 4
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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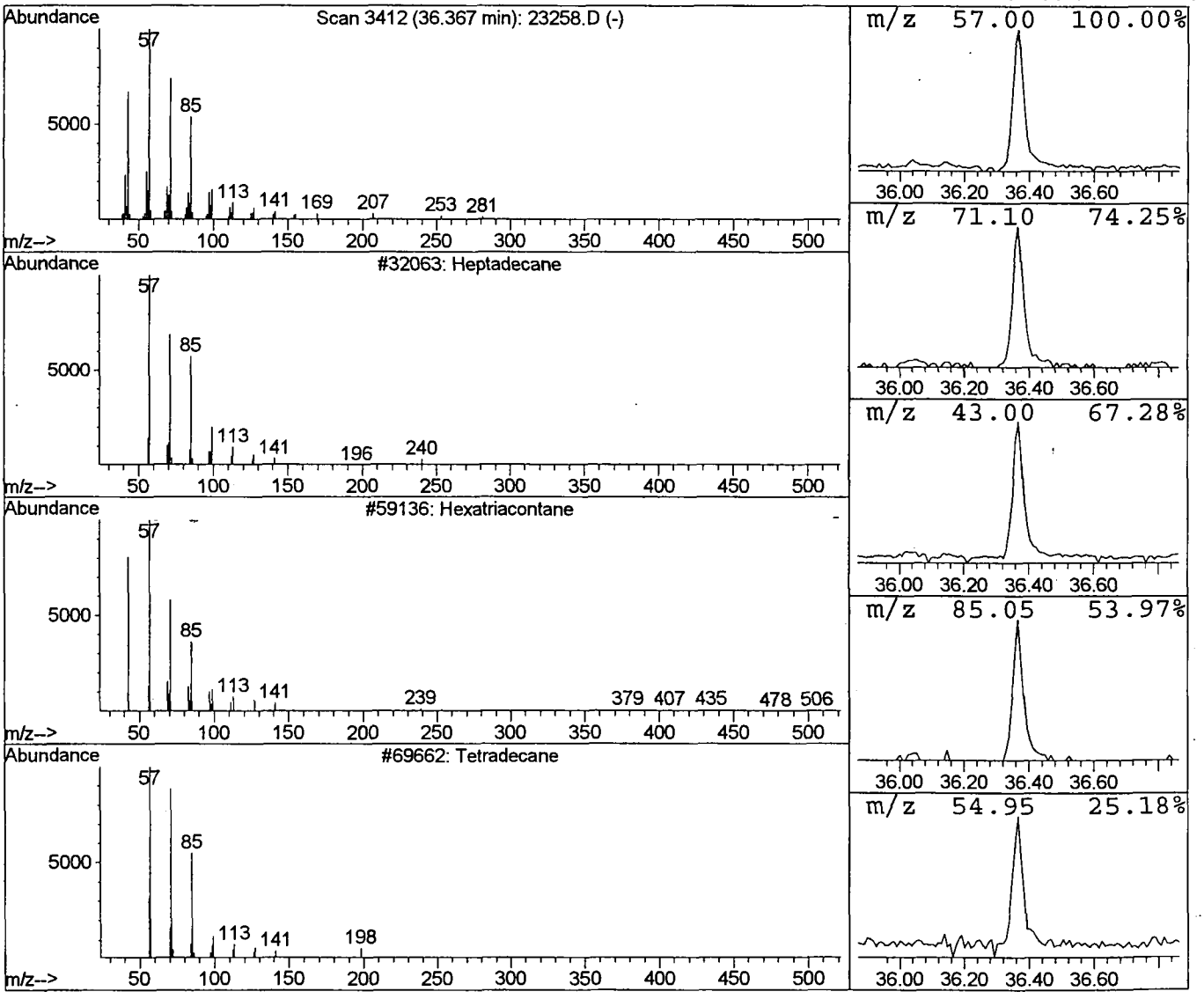
Peak Number 12 Heptadecane — Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
36.37	0.81 ug/l	123851	Perylene-d12	37.25

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heptadecane	240	C17H36	000629-78-7	91
2	Hexatriacontane	507	C36H74	000630-06-8	91
3	Tetradecane	198	C14H30	000629-59-4	91
4	Tetratetracontane	619	C44H90	007098-22-8	91



Tentatively Identified Compound (LSC) summary

Operator ID: 209 GALOS Date Acquired: 9 Oct 2001 4:17 pm
Data File: C:\HPCHEM\1\DATA\OCT0901\23258.D
Name: g c/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
2-Hexanone	6.50	0.4	ug/l	85219	ISTD01	11.76	4005080	20.0
1,4-Dichlorobenzene-	11.62	0.6	ug/l	110634	ISTD01	11.76	4005080	20.0
Butyl hexadecanoate	29.51	1.6	ug/l	295775	ISTD05	33.13	3707540	20.0
Pentacosane	29.63	0.4	ug/l	76204	ISTD05	33.13	3707540	20.0
Tetracosane	30.70	0.7	ug/l	122422	ISTD05	33.13	3707540	20.0
Octadecanoic acid, 2	31.65	1.1	ug/l	210510	ISTD05	33.13	3707540	20.0
Heneicosane	31.74	1.2	ug/l	223828	ISTD05	33.13	3707540	20.0
Heneicosane	32.73	1.3	ug/l	240258	ISTD05	33.13	3707540	20.0
Octadecane	33.69	1.5	ug/l	286357	ISTD05	33.13	3707540	20.0
Octadecane	34.61	1.2	ug/l	213473	ISTD05	33.13	3707540	20.0
Hexatriacontane	35.50	1.2	ug/l	183151	ISTD06	37.25	3044410	20.0
Heptadecane	36.37	0.8	ug/l	123851	ISTD06	37.25	3044410	20.0

23258.D NP091101.M Thu Oct 11 10:41:22 2001

Comments for Sample AD23258 - Analysis $GCMS

Westchester Dept. of Labs & Research
User: Vinci, David L
Date: 10-12-2001 Time: 17:59:24

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.