User: Galos, Marie
Westchester Dept. of Labs & Research
Date: 10/12/2001 Time: 09:22:06

Sample: AD23260
Analysis: $GCMS GCMS Scan For Unknowns
Units: ug
Started: 10/12/01 09:11 Ended: 10/12/01 09:11 Analyst: MG
Page: 1

	Component Name	Viol	Result
1	Other unkown compounds		see comment

Quantitation Report (QT Reviewed)

Data File : C:\HPCHEM\1\DATA\OCT0901\23260.D Vial: 6
Acq On : 9 Oct 2001 6:15 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 19:03 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	758707	20.00	ug/l	-0.15
19) Naphthalene-d8	15.38	136	3042315	20.00	ug/l	-0.15
31) Acenaphthene-d10	20.67	164	1430494	20.00	ug/l	-0.15
49) Phenanthrene-d10	25.09	188	2039296	20.00	ug/l	-0.15
59) Chrysene-d12	33.13	240	1223490	20.00	ug/l	-0.15
68) Perylene-d12	37.24	264	809540	20.00	ug/l	-0.18
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	11.77	132	482	0.01	ug/l	0.36
Spiked Amount 75.000			Recovery =	0.01%		
7) 1,2-Dichlorobenzene-d4	12.28	152	8939	0.24	ug/l	-0.15
Spiked Amount 50.000			Recovery =	0.48%		
20) Nitrobenzene-d5	0.00	82	0	0.00	ug/l	
Spiked Amount 50.000			Recovery =	0.00%		
32) 2-Fluorobiphenyl	18.80	172	3169	0.03	ug/l	0.00
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.		

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

Quantitation Report (QT Reviewed)

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Acq On : 9 Oct 2001 6:15 pm Operator: 209 GALOS
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Misc : Multiplr: 1.00
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
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	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	12.98	77	819	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	15.42	128	197	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	21.08	165	178	N.D.	
45)	Diethylphthalate	22.16	149	268	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	25.15	178	216	N.D.	
56)	Anthracene	25.10	178	784	N.D.	
57)	Di-n-butylphthalate	27.08	149	4544	N.D.	
58)	Fluoranthene	0.00	202	0	N.D.	
60)	3,3'-Dichlorobenzidine	0.00	252	0	N.D.	

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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
61)	Benzidine	0.00	184	0	N.D.	
63)	Pyrene	0.00	202	0	N.D.	
64)	Butylbenzylphthalate	31.60	149	727	N.D.	
65)	Benzo(a)anthracene	33.13	228	2740	N.D.	
66)	Bis(2-ethylhexyl)phthalate	33.39	149	4983	N.D.	
67)	Chrysene	33.13	228	2740	N.D.	
69)	Di-n-Octylphthalate	35.14	149	635	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	37.24	252	2652	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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Title : 209 625/TCLP calibration table 7/16/01
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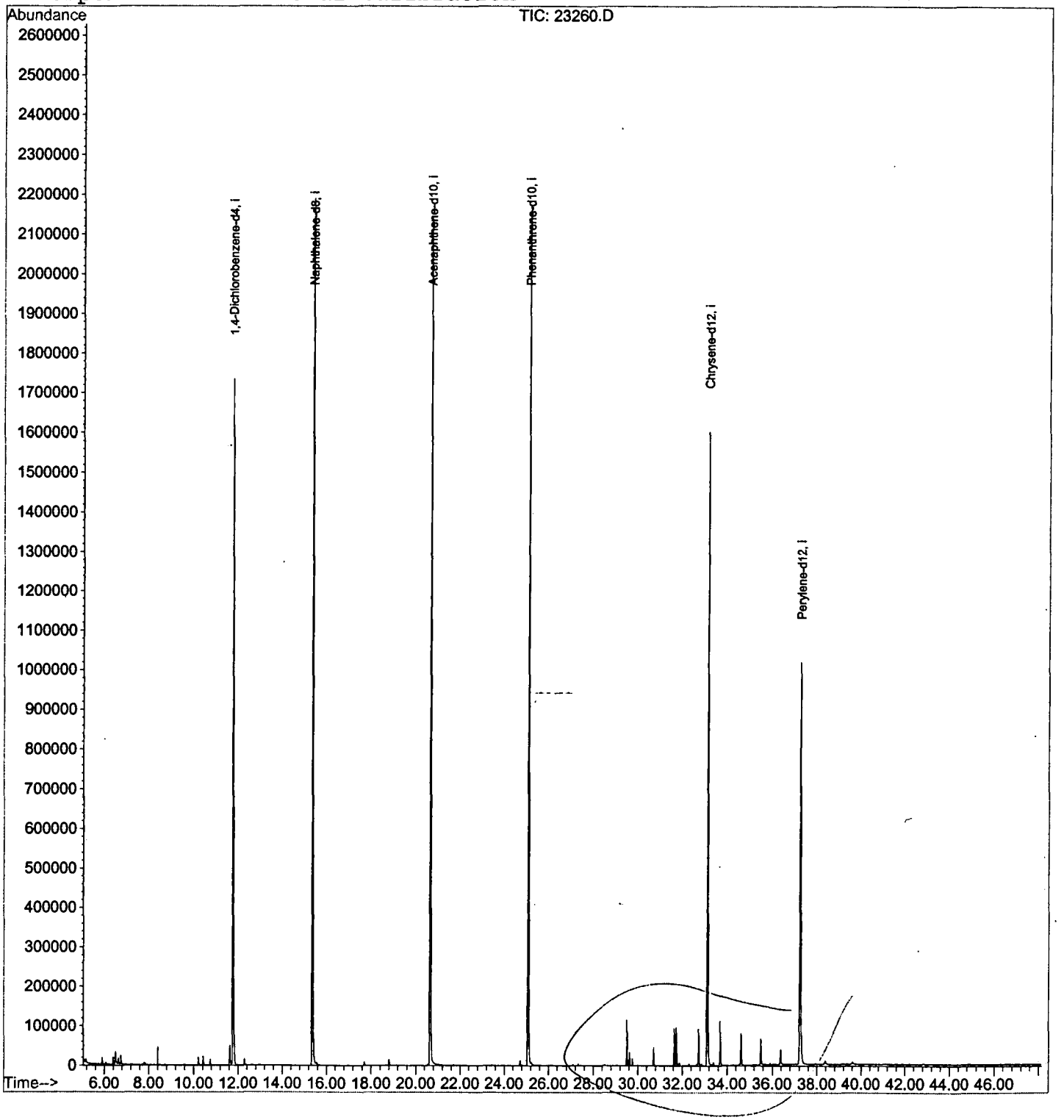
Response via : Initial Calibration



LSC Area Percent Report

Data File : C:\HPCHEM\1\DATA\OCT0901\23260.D Vial: 6
Acq On : 9 Oct 2001 6:15 pm Operator: 209 GALOS
Sample : g c/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

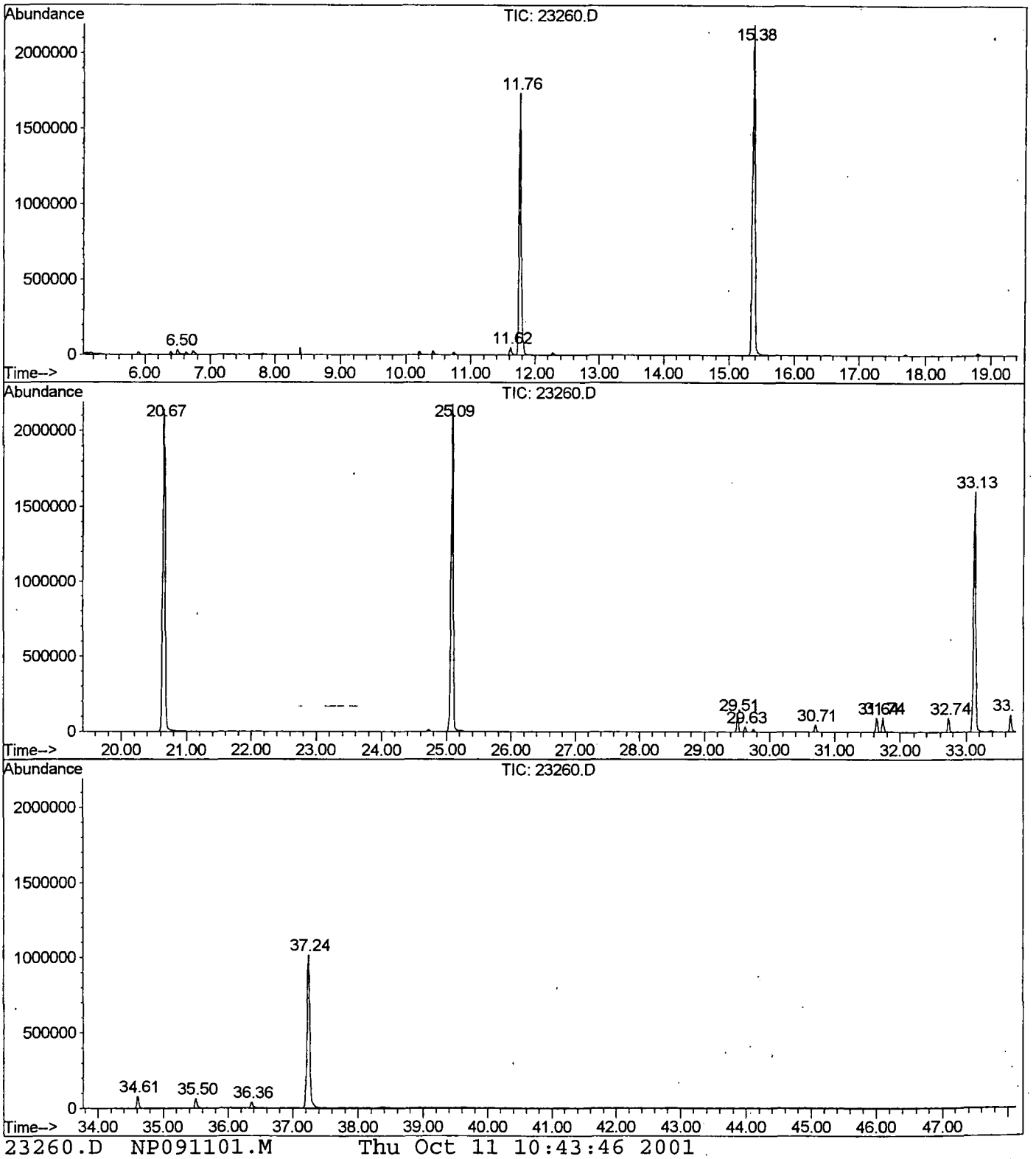
peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	peak area	peak % max.	% of total
1	6.500	154	159	170	rVV	32190	77804	1.36%	0.267%
2	11.623	713	717	726	rBV	49698	110116	1.92%	0.377%
3	11.760	726	732	750	rVV	1733710	3966951	69.28%	13.593%
4	15.377	1118	1126	1142	rBV	2189224	5576154	97.38%	19.108%
5	20.665	1693	1702	1721	rBV	2139989	5726158	100.00%	19.622%
6	25.090	2174	2184	2196	rBV2	2173375	5240062	91.51%	17.956%
7	29.506	2659	2665	2673	rBV	114463	231258	4.04%	0.792%
8	29.625	2673	2678	2684	rVV	33373	62058	1.08%	0.213%
9	30.709	2788	2796	2802	rBV	46500	94234	1.65%	0.323%
10	31.645	2892	2898	2903	rBV	92604	174776	3.05%	0.599%
11	31.737	2903	2908	2914	rVV	94697	184067	3.21%	0.631%
12	32.737	3008	3017	3023	rBV	92512	198876	3.47%	0.681%
13	33.132	3050	3060	3074	rBV2	1601818	3831720	66.92%	13.130%
14	33.692	3111	3121	3132	rBV	112099	242454	4.23%	0.831%
15	34.610	3217	3221	3229	rVB	78097	166335	2.90%	0.570%
16	35.501	3310	3318	3327	rBV	66176	158129	2.76%	0.542%
17	36.364	3405	3412	3421	rBV	37813	95551	1.67%	0.327%
18	37.245	3498	3508	3528	rBV2	1014989	3046225	53.20%	10.438%

Sum of corrected areas: 29182928

23260.D NP091101.M Thu Oct 11 10:43:43 2001

LSC Report - Integrated Chromatogram

File : C:\HPCHEM\1\DATA\OCT0901\23260.D
Operator : 209 GALOS
Acquired : 9 Oct 2001 6:15 pm using AcqMethod NP091101
Instrument : GC/MS 209
Sample Name: g c/ms unknown
Misc Info :
Vial Number: 6
Quant File :NP091101.RES (RTE Integrator)



23260.D NP091101.M Thu Oct 11 10:43:46 2001

Library Search Compound Report

Data File : C:\HPCHEM\1\DATA\OCT0901\23260.D
Acq On : 9 Oct 2001 6:15 pm
Sample : g c/ms unknown
Misc :
MS Integration Params: LSCINT.P

Vial: 6
Operator: 209 GALOS
Inst : GC/MS 209
Multiplr: 1.00

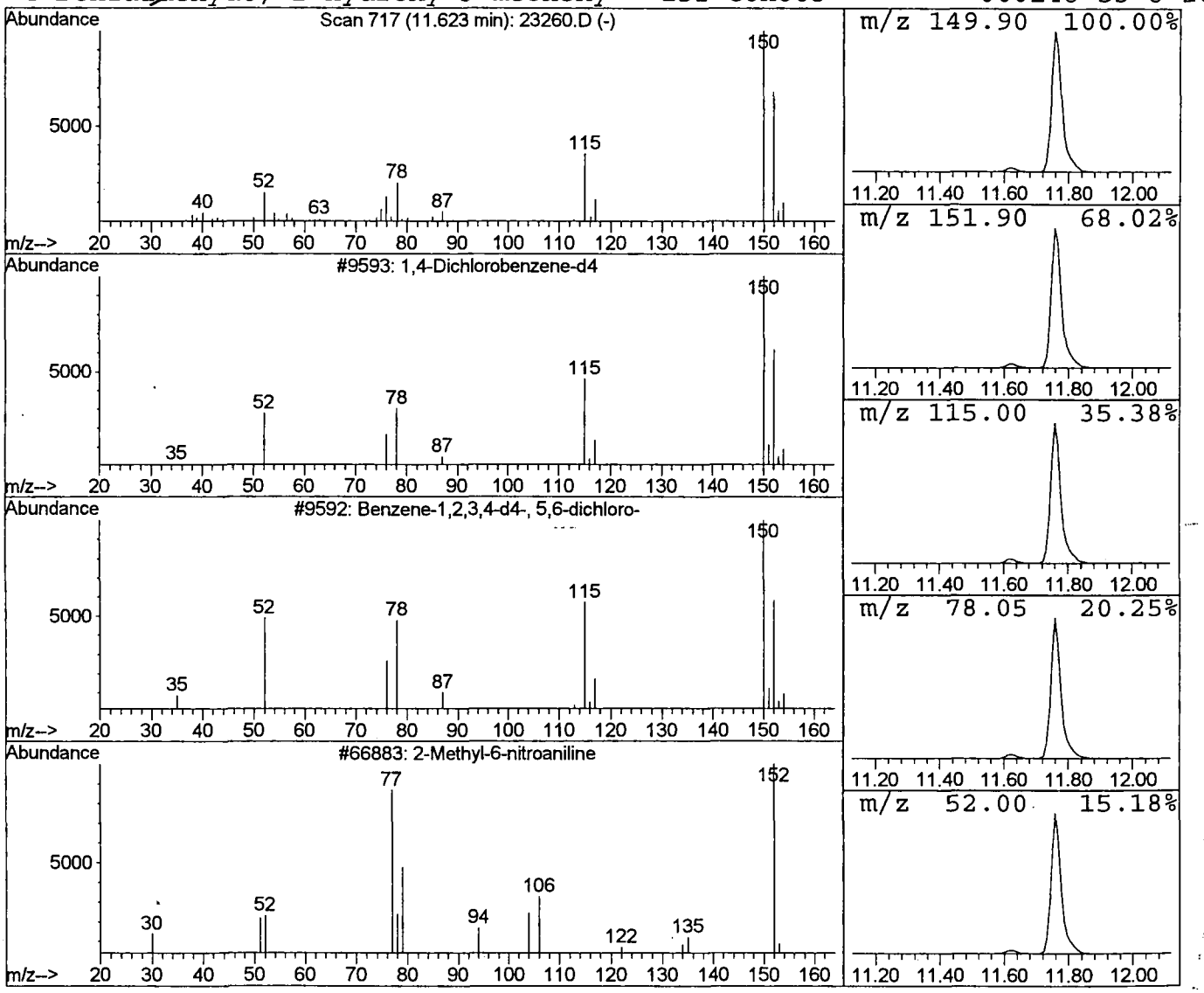
Quant Method : C:\HPCHEM\1\METHODS\NP091101.M (RTE Integrator)
Title : 209 625/TCLP calibration table 7/16/01
Library : C:\DATABASE\NBS75K.L

****

Peak Number 1 1,4-Dichlorobenzene-d4 Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.62	0.56 ug/l	110116	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	35
3	2-Methyl-6-nitroaniline	152	C7H8N2O2	000570-24-1	12
4	Benzaldehyde, 2-hydroxy-3-methoxy-	152	C8H8O3	000148-53-8	10



Library Search Compound Report

Data File : C:\HPCHEM\1\DATA\OCT0901\23260.D
Acq On : 9 Oct 2001 6:15 pm
Sample : g c/ms unknown
Misc :
MS Integration Params: LSCINT.P

Vial: 6
Operator: 209 GALOS
Inst : GC/MS 209
Multiplr: 1.00

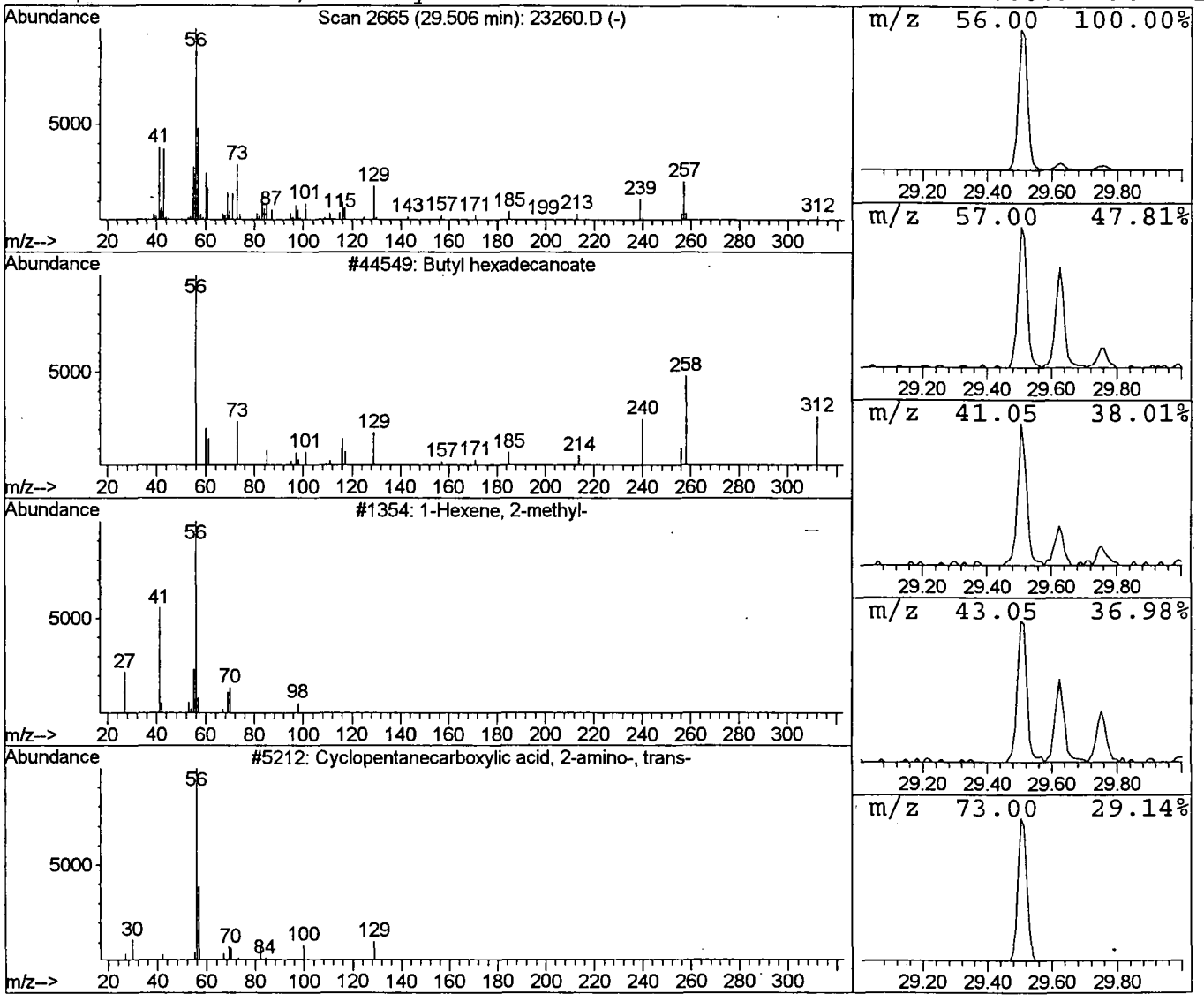
Quant Method : C:\HPCHEM\1\METHODS\NP091101.M (RTE Integrator)
Title : 209 625/TCLP calibration table 7/16/01
Library : C:\DATABASE\NBS75K.L

****

Peak Number 2 Butyl hexadecanoate Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
29.51	1.21 ug/l	231258	Chrysene-d12	33.13

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Butyl hexadecanoate	312	C20H40O2	000000-00-0	35
2	1-Hexene, 2-methyl-	98	C7H14	006094-02-6	27
3	Cyclopentanecarboxylic acid, 2-amin	129	C6H11NO2	040482-05-1	23
4	1,3-Hexanediol, 2-ethyl-	146	C8H18O2	000094-96-2	23



Library Search Compound Report

Data File : C:\HPCHEM\1\DATA\OCT0901\23260.D
Acq On : 9 Oct 2001 6:15 pm
Sample : g c/ms unknown
Misc :
MS Integration Params: LSCINT.P

Vial: 6
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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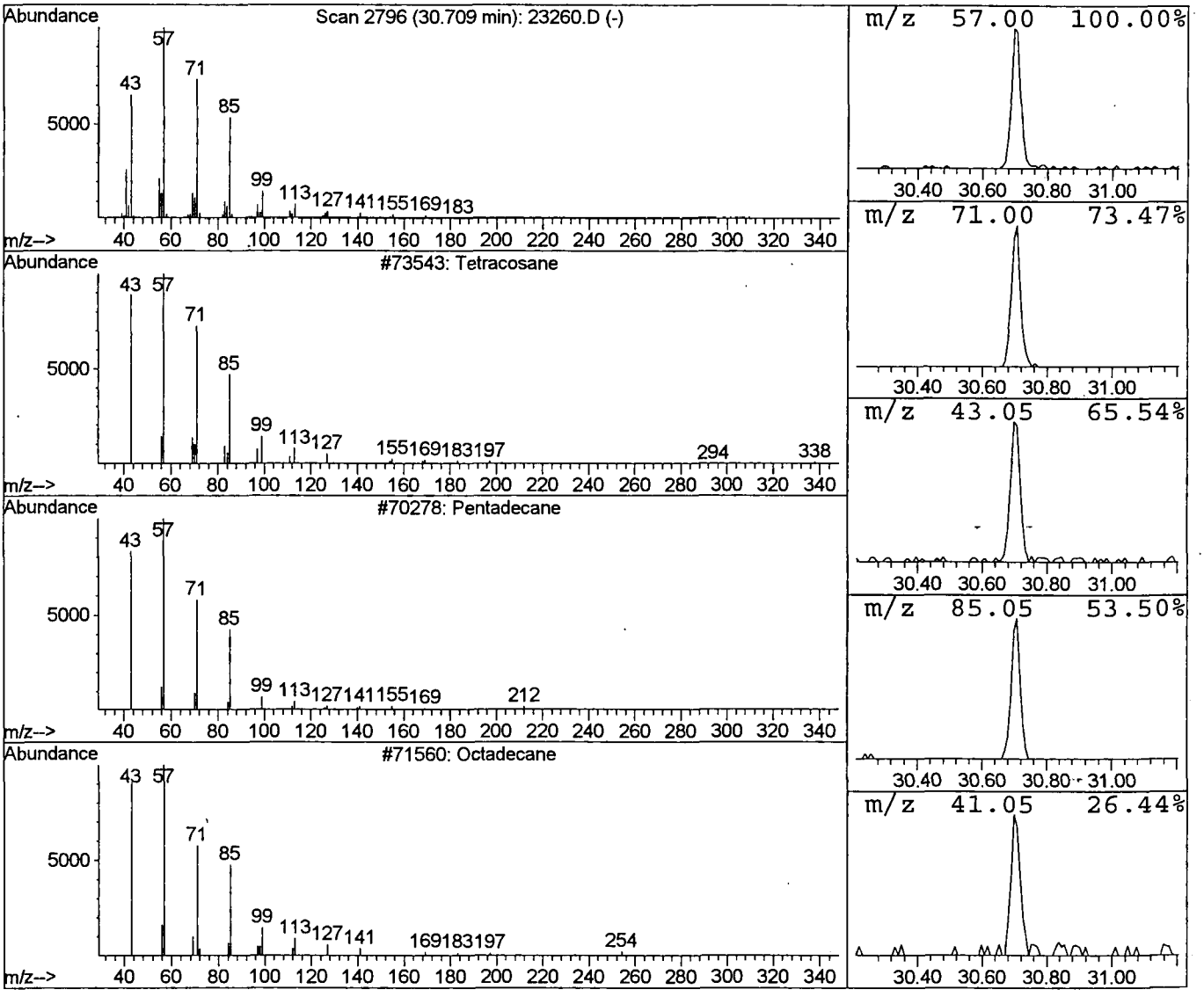
Peak Number 3 Tetracosane — Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
30.71	0.49 ug/l	94234	Chrysene-d12	33.13

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Tetracosane	338	C24H50	000646-31-1	90
2	Pentadecane	212	C15H32	000629-62-9	90
3	Octadecane	254	C18H38	000593-45-3	87
4	Pentacosane	352	C25H52	000629-99-2	87



Library Search Compound Report

Data File : C:\HPCHEM\1\DATA\OCT0901\23260.D
Acq On : 9 Oct 2001 6:15 pm
Sample : g c/ms unknown
Misc :
MS Integration Params: LSCINT.P

Vial: 6
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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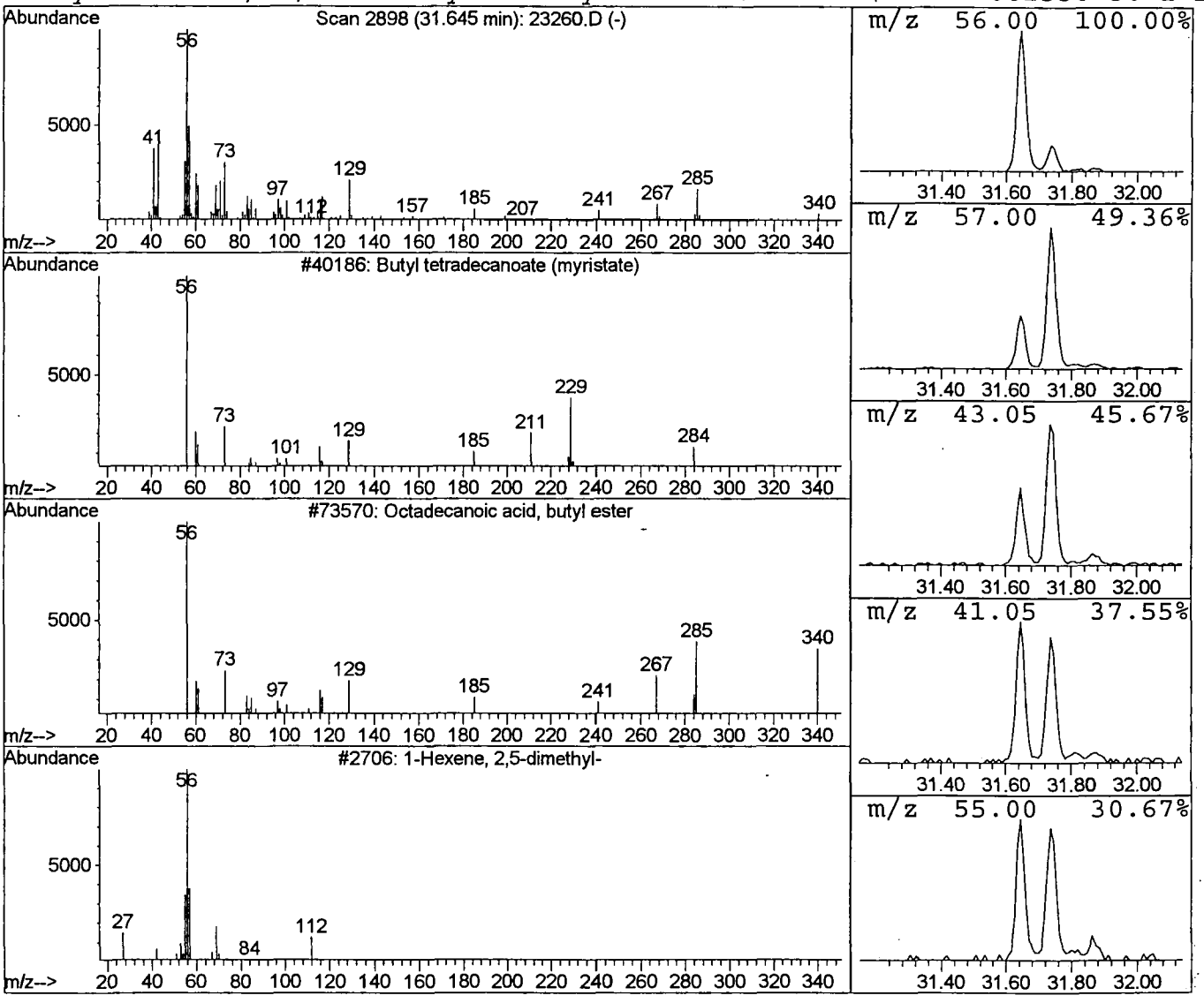
Peak Number 4 Butyl tetradecanoate (myristat Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
31.65	0.91 ug/l	174776	Chrysene-d12	33.13

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Butyl tetradecanoate (myristate)	284	C18H36O2	000000-00-0	80
2	Octadecanoic acid, butyl ester	340	C22H44O2	000123-95-5	28
3	1-Hexene, 2,5-dimethyl-	112	C8H16	006975-92-4	27
4	Cyclobutane, 1,1-dimethyl-2-octyl-	196	C14H28	062338-30-1	25



Library Search Compound Report

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Acq On : 9 Oct 2001 6:15 pm
Sample : g c/ms unknown
Misc :
MS Integration Params: LSCINT.P

Vial: 6
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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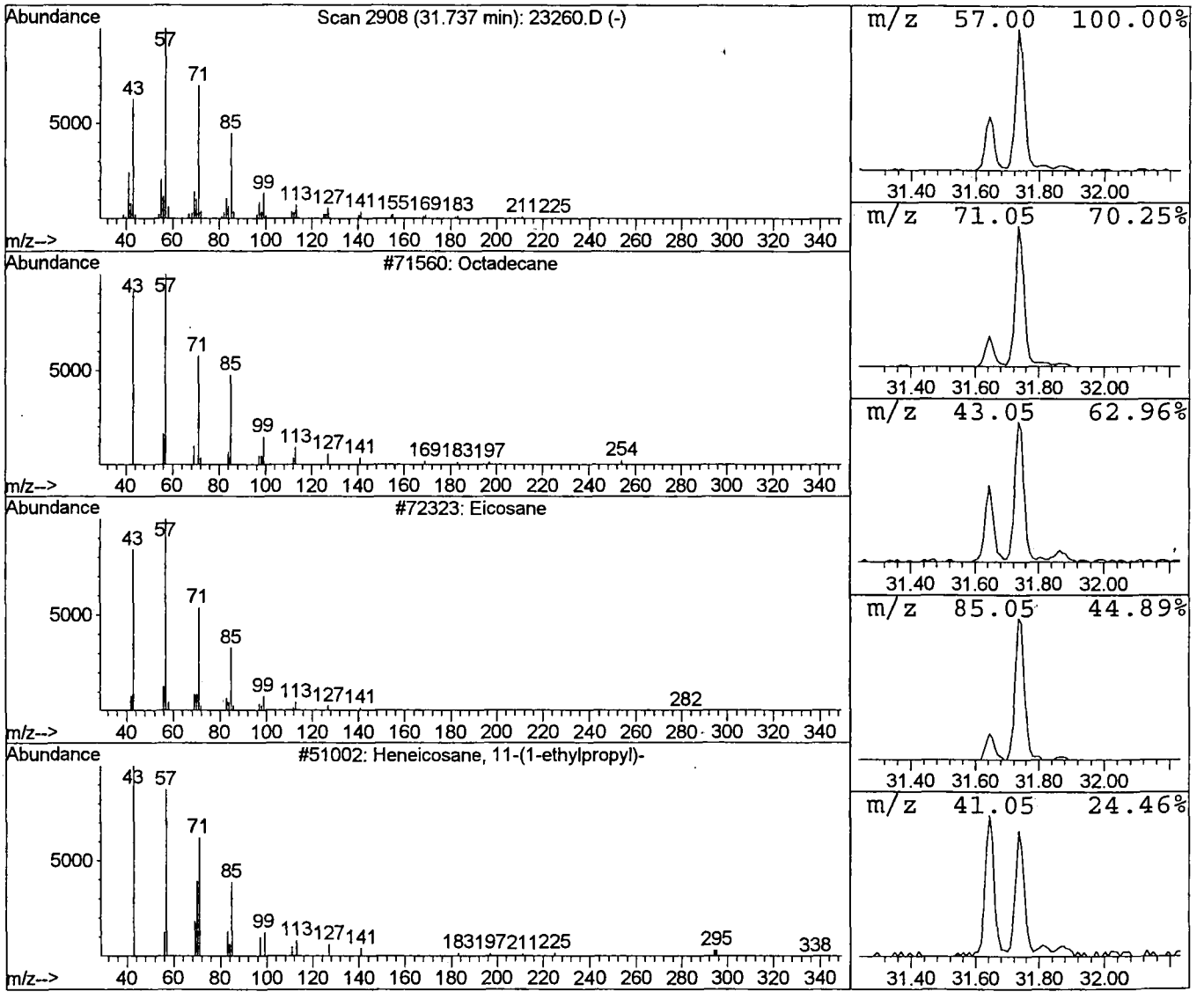
Peak Number 5 Octadecane — Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
31.74	0.96 ug/l	184067	Chrysene-d12	33.13

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Octadecane	254	C18H38	000593-45-3	91
2	Eicosane	282	C20H42	000112-95-8	91
3	Heneicosane, 11-(1-ethylpropyl)-	366	C26H54	055282-11-6	91
4	Tetracosane	338	C24H50	000646-31-1	90



Library Search Compound Report

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

Vial: 6
Operator: 209 GALOS
Inst : GC/MS 209
Multiplr: 1.00

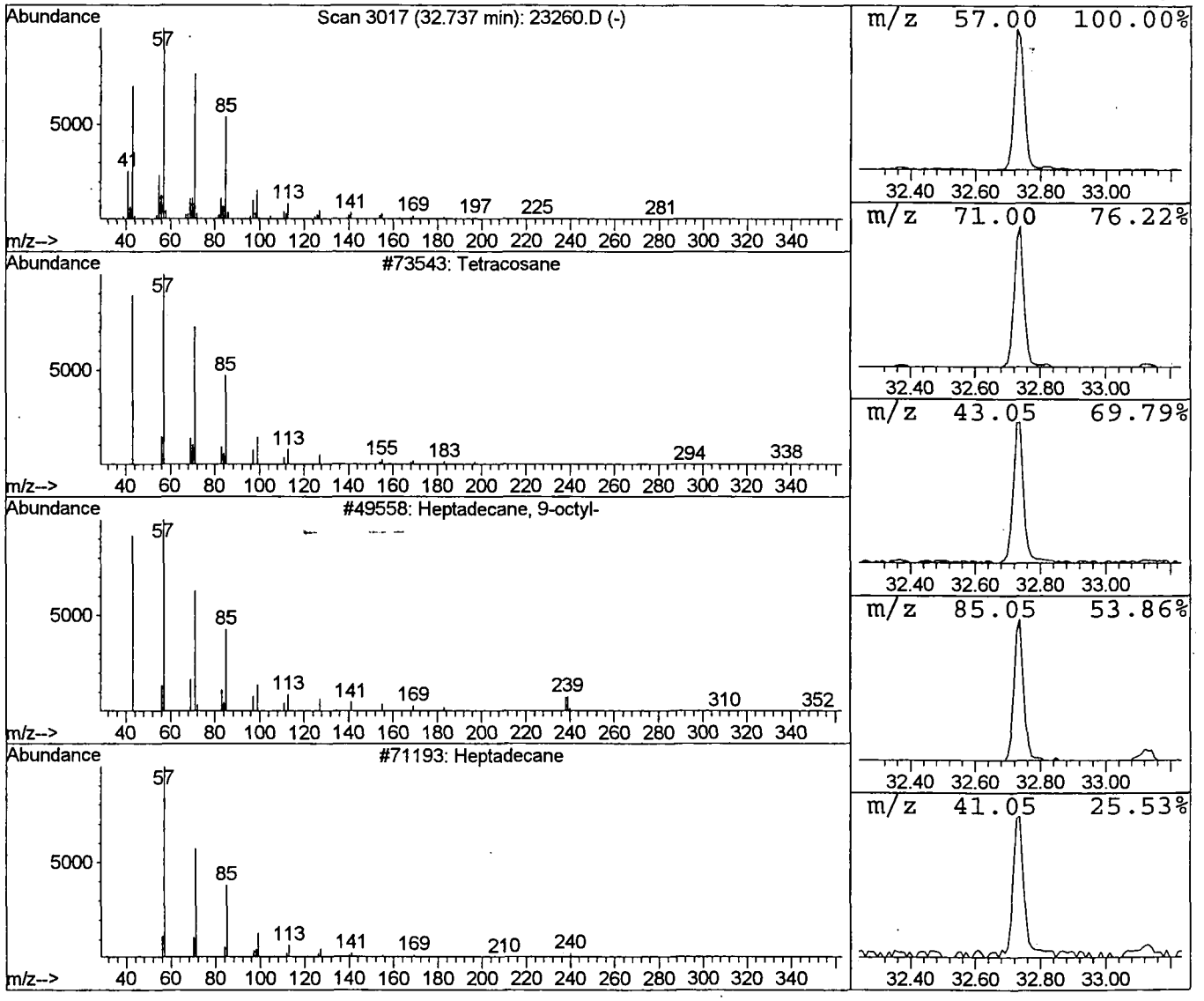
Quant Method : C:\HPCHEM\1\METHODS\NP091101.M (RTE Integrator)
Title : 209 625/TCLP calibration table 7/16/01
Library : C:\DATABASE\NBS75K.L

****

Peak Number 6 Tetracosane Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
32.74	1.04 ug/l	198876	Chrysene-d12	33.13

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Tetracosane	338	C24H50	000646-31-1	91
2	Heptadecane, 9-octyl-	352	C25H52	007225-64-1	91
3	Heptadecane	240	C17H36	000629-78-7	91
4	Pentadecane	212	C15H32	000629-62-9	86



Library Search Compound Report

Data File : C:\HPCHEM\1\DATA\OCT0901\23260.D
Acq On : 9 Oct 2001 6:15 pm
Sample : g c/ms unknown
Misc :
MS Integration Params: LSCINT.P

Vial: 6
Operator: 209 GALOS
Inst : GC/MS 209
Multiplr: 1.00

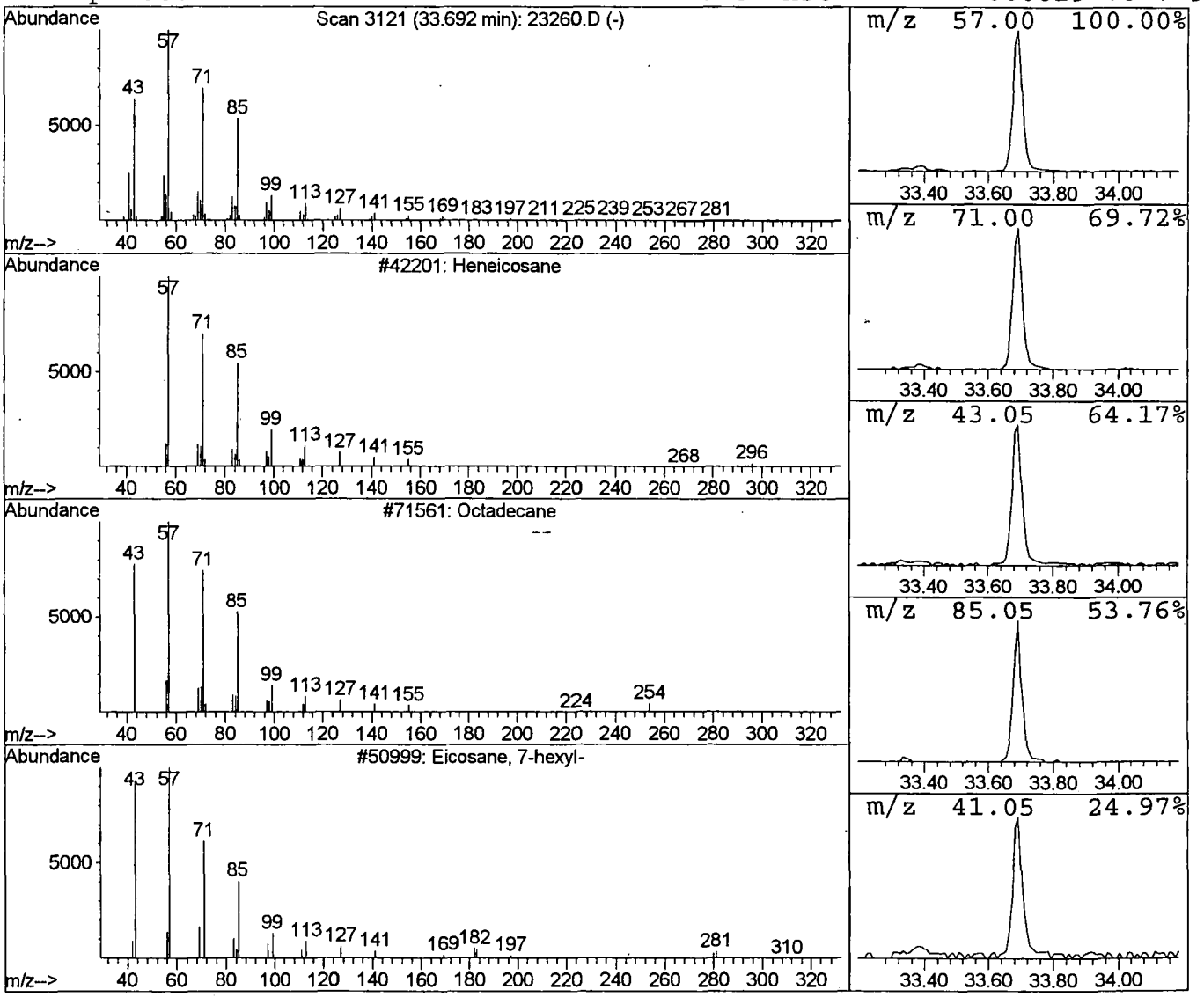
Quant Method : C:\HPCHEM\1\METHODS\NP091101.M (RTE Integrator)
Title : 209 625/TCLP calibration table 7/16/01
Library : C:\DATABASE\NBS75K.L

****

Peak Number 7 Heneicosane — Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
33.69	1.27 ug/l	242454	Chrysene-d12	33.13

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heneicosane	296	C21H44	000629-94-7	91
2	Octadecane	254	C18H38	000593-45-3	91
3	Eicosane, 7-hexyl-	366	C26H54	055333-99-8	91
4	Heptadecane	240	C17H36	000629-78-7	91



Library Search Compound Report

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Acq On : 9 Oct 2001 6:15 pm
Sample : g c/ms unknown
Misc :
MS Integration Params: LSCINT.P

Vial: 6
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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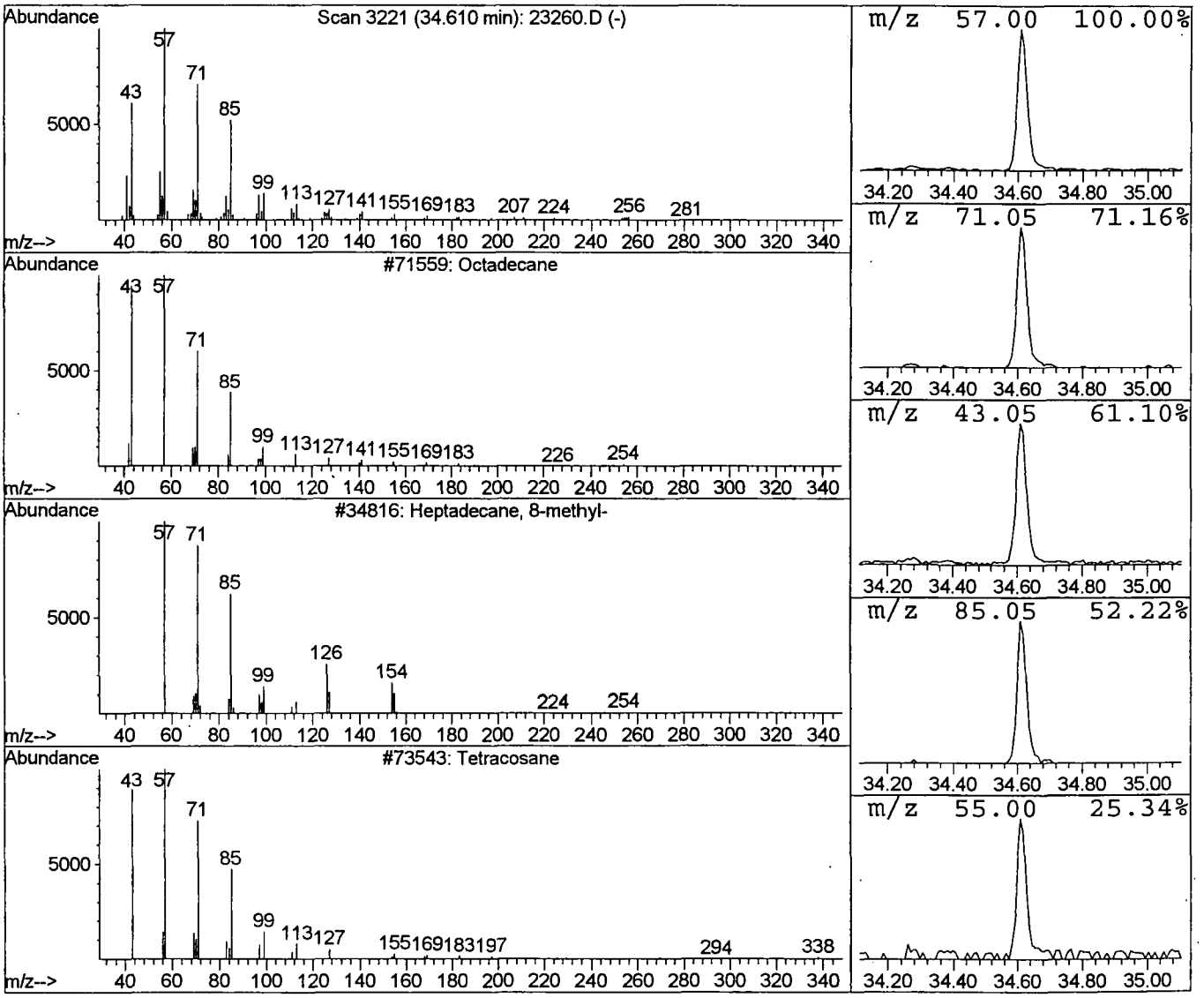
Peak Number 8 Octadecane — Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
34.61	0.87 ug/l	166335	Chrysene-d12	33.13

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Octadecane	254	C18H38	000593-45-3	93
2	Heptadecane, 8-methyl-	254	C18H38	013287-23-5	91
3	Tetracosane	338	C24H50	000646-31-1	90
4	Heneicosane	296	C21H44	000629-94-7	90



Library Search Compound Report

Data File : C:\HPCHEM\1\DATA\OCT0901\23260.D
Acq On : 9 Oct 2001 6:15 pm
Sample : g c/ms unknown
Misc :
MS Integration Params: LSCINT.P

Vial: 6
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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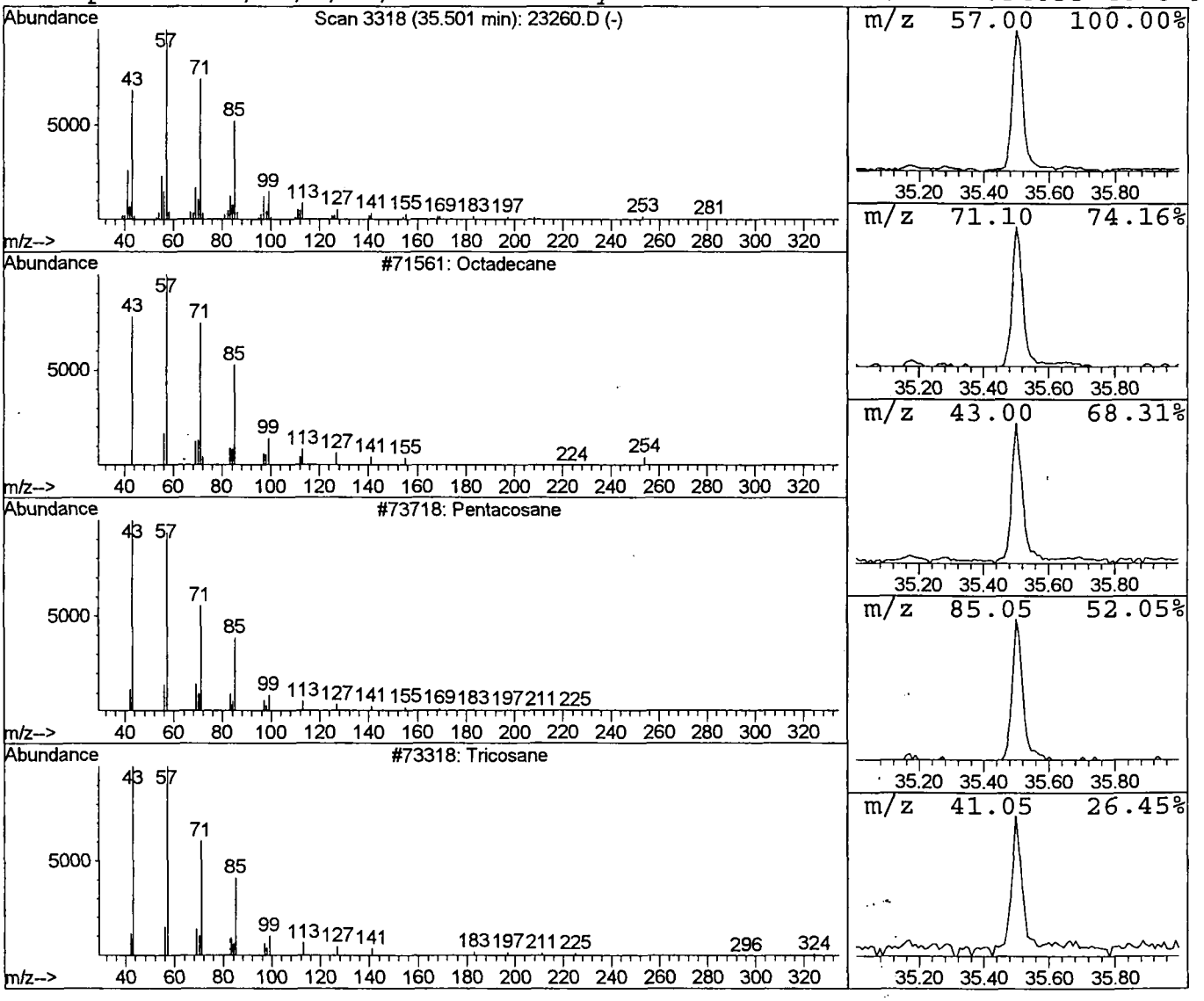
Peak Number 9 Octadecane — Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
35.50	1.04 ug/l	158129	Perylene-d12	37.24

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Octadecane	254	C18H38	000593-45-3	91
2	Pentacosane	352	C25H52	000629-99-2	91
3	Tricosane	324	C23H48	000638-67-5	91
4	Heptadecane, 2,6,10,15-tetramethyl-	296	C21H44	054833-48-6	91



Library Search Compound Report

Data File : C:\HPCHEM\1\DATA\OCT0901\23260.D
Acq On : 9 Oct 2001 6:15 pm
Sample : g c/ms unknown
Misc :
MS Integration Params: LSCINT.P

Vial: 6
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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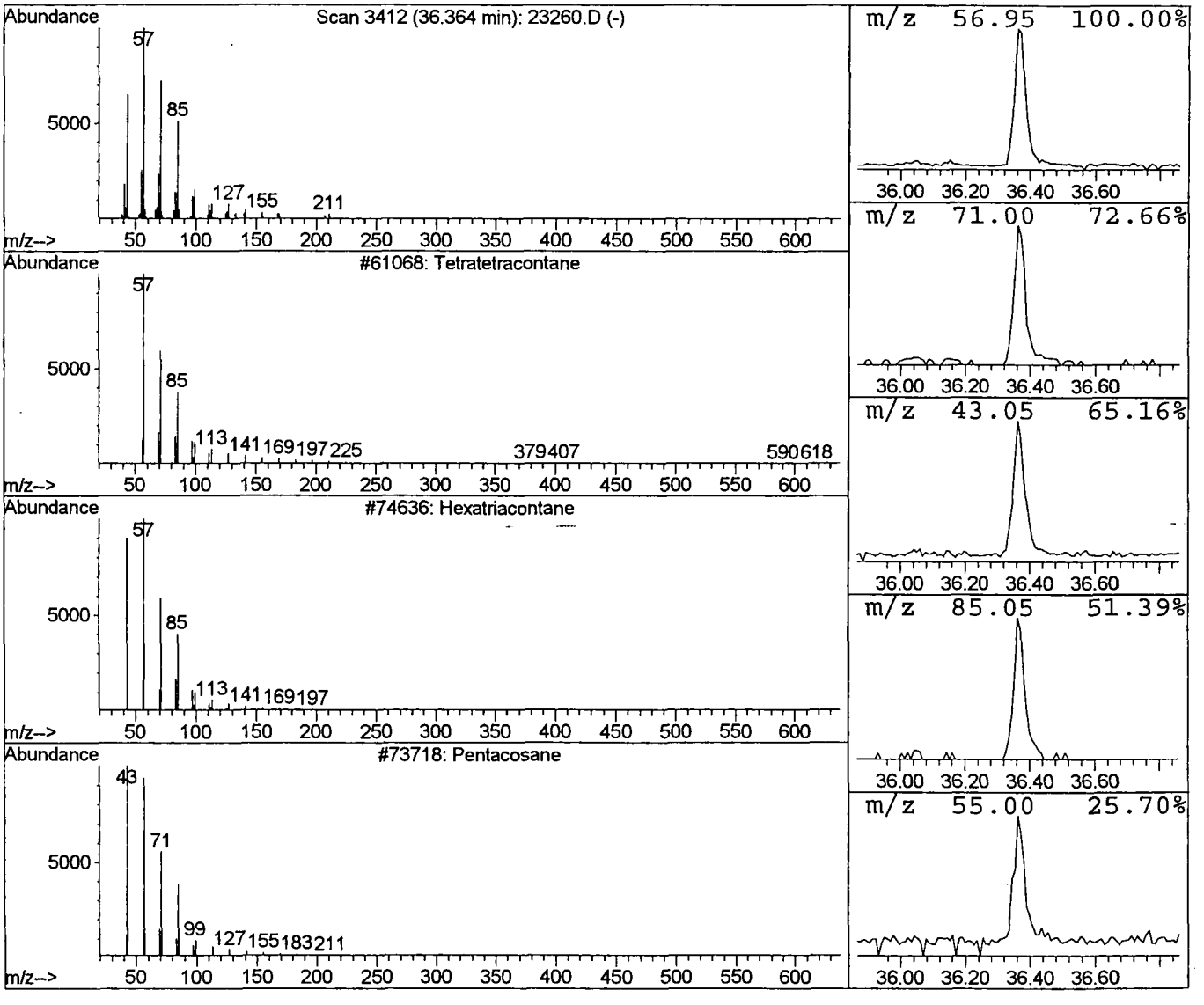
Peak Number 10 Tetratetracontane Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
36.36	0.63 ug/l	95551	Perylene-d12	37.24

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Tetratetracontane	619	C44H90	007098-22-8	91
2	Hexatriacontane	507	C36H74	000630-06-8	91
3	Pentacosane	352	C25H52	000629-99-2	91
4	Eicosane	282	C20H42	000112-95-8	91



Tentatively Identified Compound (LSC) summary

Operator ID: 209 GALOS Date Acquired: 9 Oct 2001 6:15 pm
Data File: C:\HPCHEM\1\DATA\OCT0901\23260.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
1,4-Dichlorobenzene-	11.62	0.6	ug/l	110116	ISTD01	11.76	3966950	20.0
Butyl hexadecanoate	29.51	1.2	ug/l	231258	ISTD05	33.13	3831720	20.0
Tetracosane	30.71	0.5	ug/l	94234	ISTD05	33.13	3831720	20.0
Butyl tetradecanoate	31.65	0.9	ug/l	174776	ISTD05	33.13	3831720	20.0
Octadecane	31.74	1.0	ug/l	184067	ISTD05	33.13	3831720	20.0
Tetracosane	32.74	1.0	ug/l	198876	ISTD05	33.13	3831720	20.0
Heneicosane	33.69	1.3	ug/l	242454	ISTD05	33.13	3831720	20.0
Octadecane	34.61	0.9	ug/l	166335	ISTD05	33.13	3831720	20.0
Octadecane	35.50	1.0	ug/l	158129	ISTD06	37.24	3046230	20.0
Tetratetracontane	36.36	0.6	ug/l	95551	ISTD06	37.24	3046230	20.0

23260.D NP091101.M Thu Oct 11 10:44:10 2001

Comments for Sample AD23260 - Analysis $GCMS

Westchester Dept. of Labs & Research
User: Vinci, David L
Date: 10-12-2001 Time: 18:04:45

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.