User: Galos, Marie
Westchester Dept. of Labs & Research
Date: 10/16/2001 Time: 09:21:55

Sample: AD23343
Analysis: $GCMS GCMS Scan For Unknowns
Units: ug
Started: 10/16/01 09:21 Ended: 10/16/01 09:21 Analyst: MG
Page: 1

	Component Name	Viol	Result
1	Other unknown compounds		see comment

analysis 10/10 - bad blanks flew.

Quantitation Report (QT Reviewed)

Data File : C:\HPCHEM\1\DATA\OCT0901\23343.D Vial: 16
Acq On : 10 Oct 2001 3:59 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 10:30 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	375982	20.00	ug/l	-0.15
19) Naphthalene-d8	15.38	136	1489588	20.00	ug/l	-0.15
31) Acenaphthene-d10	20.65	164	686125	20.00	ug/l	-0.16
49) Phenanthrene-d10	25.09	188	941024	20.00	ug/l	-0.15
59) Chrysene-d12	33.12	240	595185	20.00	ug/l	-0.16
68) Perylene-d12	37.23	264	440116	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	0.00	132	0	0.00	ug/l	
Spiked Amount 75.000			Recovery =	0.00%		
7) 1,2-Dichlorobenzene-d4	12.28	152	3982	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	18.79	172	1358	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	5.12	79	180	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	12.97	77	485	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
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	22.16	149	182	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	2472	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	1283	N.D.	
66)	Bis(2-ethylhexyl)phthalate	33.39	149	2132	N.D.	
67)	Chrysene	33.12	228	1283	N.D.	
69)	Di-n-Octylphthalate	0.00	149	0	N.D.	
70)	Benzo(b)fluoranthene	0.00	252	0	N.D.	

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

Quantitation Report (QT Reviewed)

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Sample : gc/ms unknown Inst : GC/MS 209
Misc : Multiplr: 1.00
MS Integration Params: RTEINT.P
Quant Time: Oct 11 10:30 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	0.00	252	0	N.D.	
72)	Benzo(a)pyrene	37.24	252	1461	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.	

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

Quantitation Report

Data File : C:\HPCHEM\1\DATA\OCT0901\23343.D Vial: 16
Acq On : 10 Oct 2001 3:59 am Operator: 209 GALOS
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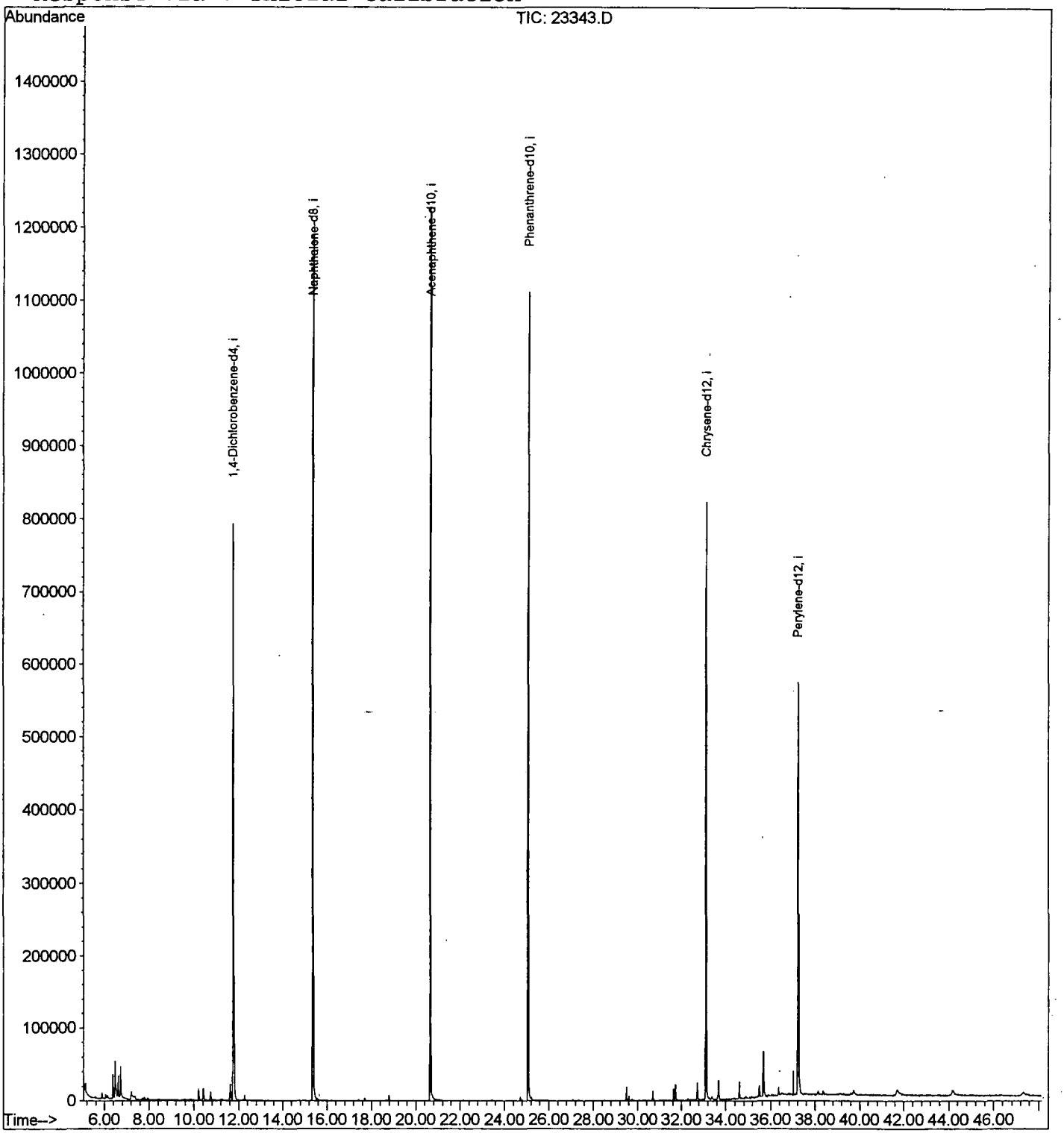
Title : 209 625/TCLP calibration table 7/16/01
Last Update : Thu Sep 13 12:47:04 2001
Response via : Initial Calibration



LSC Area Percent Report

Data File : C:\HPCHEM\1\DATA\OCT0901\23343.D Vial: 16
Acq On : 10 Oct 2001 3:59 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

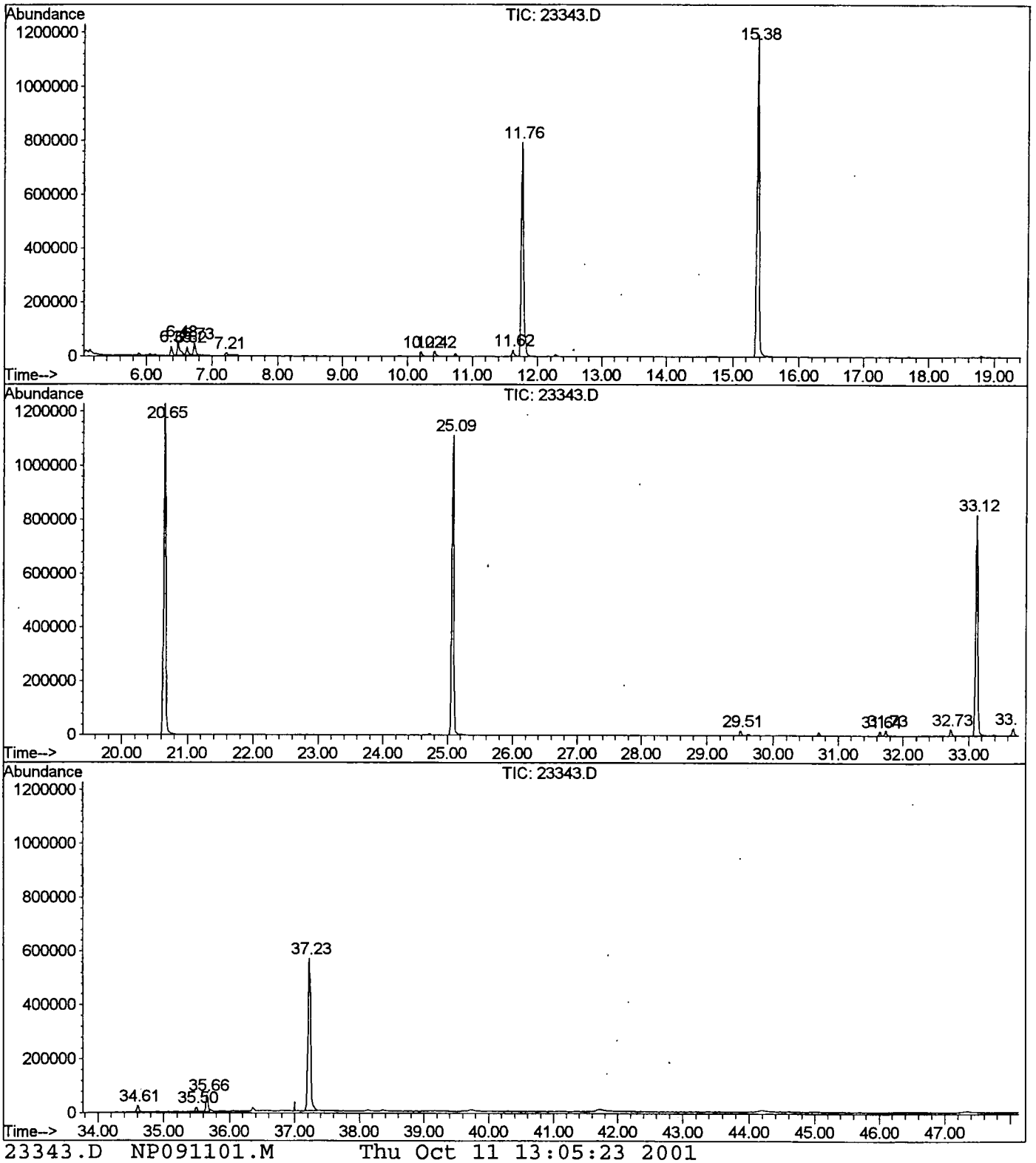
peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	peak area	peak % max.	% of total
1	6.379	142	146	152	rBV	33387	63697	2.33%	0.444%
2	6.480	154	157	168	rVV2	51372	139049	5.08%	0.969%
3	6.617	168	172	178	rVV2	30108	63357	2.32%	0.442%
4	6.727	180	184	194	rVB	42225	89838	3.28%	0.626%
5	7.214	230	237	243	rBV2	10944	34860	1.27%	0.243%
6	10.216	559	564	574	rVB	15882	35610	1.30%	0.248%
7	10.418	582	586	594	rBV	16419	35687	1.30%	0.249%
8	11.621	712	717	727	rVB	22272	53447	1.95%	0.373%
9	11.758	727	732	747	rBV	792664	1962255	71.72%	13.680%
10	15.375	1119	1126	1139	rBV	1188714	2711389	99.10%	18.903%
11	20.654	1694	1701	1714	rBV	1227959	2736116	100.00%	19.075%
12	25.088	2176	2184	2195	rBV2	1111027	2459932	89.91%	17.150%
13	29.513	2657	2666	2670	rBV	18759	38344	1.40%	0.267%
14	31.643	2890	2898	2904	rBV	15734	32816	1.20%	0.229%
15	31.735	2904	2908	2914	rVV	21269	41539	1.52%	0.290%
16	32.726	3009	3016	3023	rBV2	23395	49818	1.82%	0.347%
17	33.121	3052	3059	3069	rBV2	820879	1862630	68.08%	12.986%
18	33.690	3110	3121	3127	rBV2	26421	67583	2.47%	0.471%
19	34.608	3216	3221	3226	rVB2	22840	49254	1.80%	0.343%
20	35.499	3313	3318	3323	rBV3	16006	35735	1.31%	0.249%
21	35.664	3331	3336	3340	rBV	61294	135080	4.94%	0.942%
22	37.234	3498	3507	3522	rBV2	565833	1645629	60.14%	11.473%

Sum of corrected areas: 14343665

23343.D NP091101.M Thu Oct 11 13:05:20 2001

LSC Report - Integrated Chromatogram

File : C:\HPCHEM\1\DATA\OCT0901\23343.D
Operator : 209 GALOS
Acquired : 10 Oct 2001 3:59 am using AcqMethod NP091101
Instrument : GC/MS 209
Sample Name: gc/ms unknown
Misc Info :
Vial Number: 16
Quant File :NP091101.RES (RTE Integrator)



23343.D NP091101.M Thu Oct 11 13:05:23 2001

Library Search Compound Report

Data File : C:\HPCHEM\1\DATA\OCT0901\23343.D
Acq On : 10 Oct 2001 3:59 am
Sample : gc/ms unknown
Misc :
MS Integration Params: LSCINT.P

Vial: 16
Operator: 209 GALOS
Inst : GC/MS 209
Multiplr: 1.00

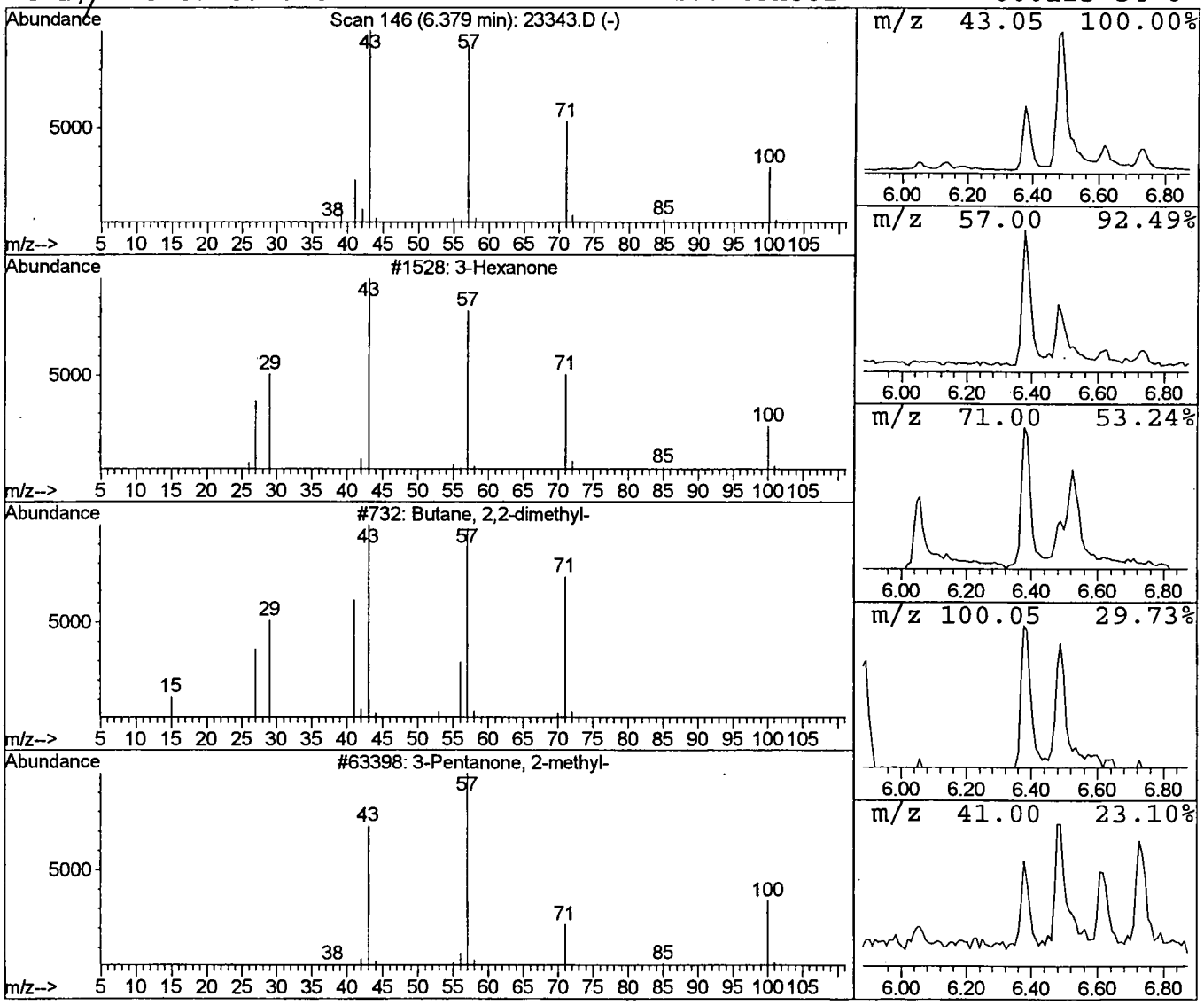
Quant Method : C:\HPCHEM\1\METHODS\NP091101.M (RTE Integrator)
Title : 209 625/TCLP calibration table 7/16/01
Library : C:\DATABASE\NBS75K.L

****

Peak Number 1 3-Hexanone Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.38	0.65 ug/l	63697	1,4-Dichlorobenzene-d4	11.76

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	3-Hexanone	100	C6H12O	000589-38-8	91
2	Butane, 2,2-dimethyl-	86	C6H14	000075-83-2	50
3	3-Pentanone, 2-methyl-	100	C6H12O	000565-69-5	47
4	2,4-Pentanedione	100	C5H8O2	000123-54-6	43



Library Search Compound Report

Data File : C:\HPCHEM\1\DATA\OCT0901\23343.D
Acq On : 10 Oct 2001 3:59 am
Sample : gc/ms unknown
Misc :
MS Integration Params: LSCINT.P

Vial: 16
Operator: 209 GALOS
Inst : GC/MS 209
Multiplr: 1.00

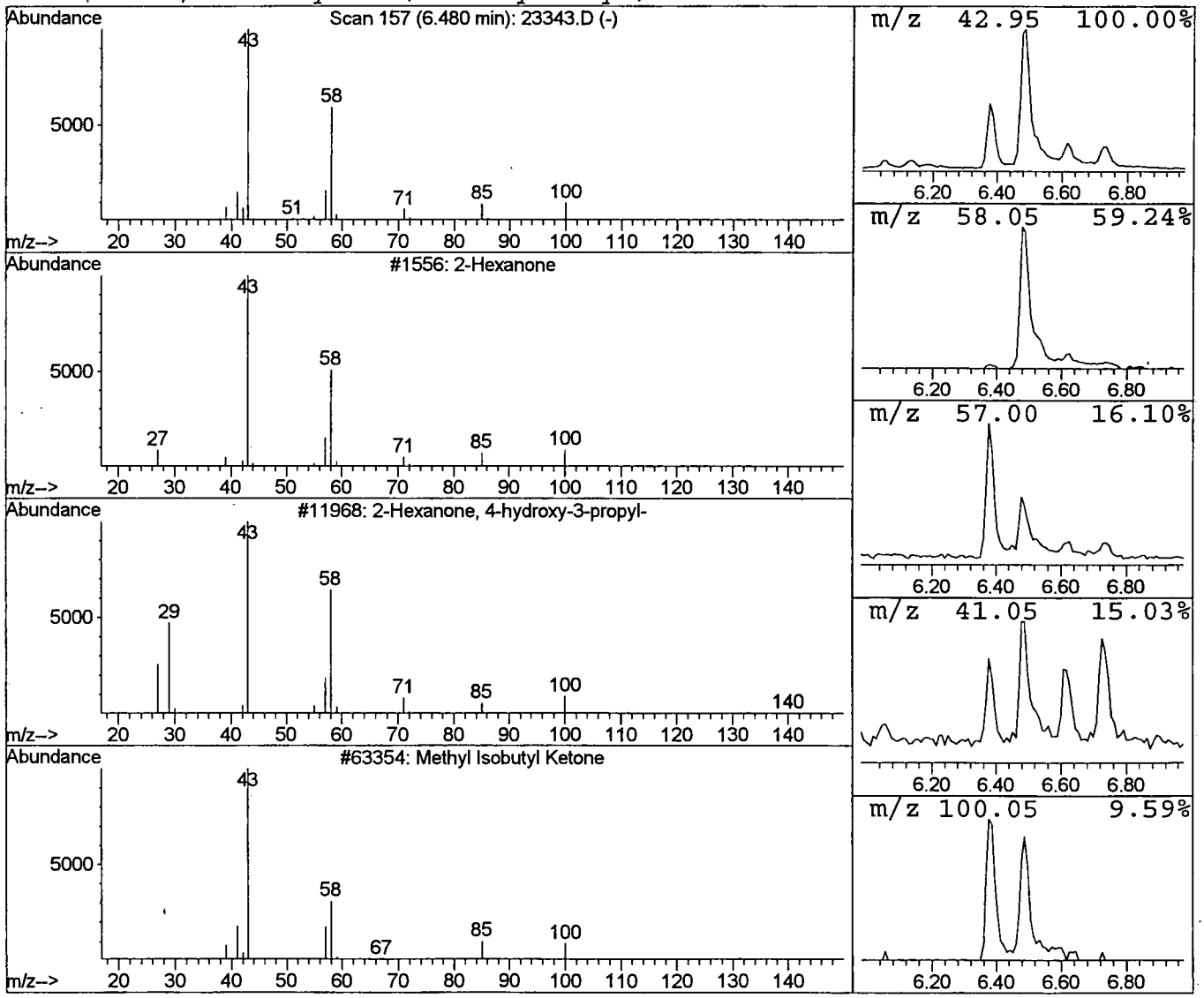
Quant Method : C:\HPCHEM\1\METHODS\NP091101.M (RTE Integrator)
Title : 209 625/TCLP calibration table 7/16/01
Library : C:\DATABASE\NBS75K.L

****

Peak Number 2 2-Hexanone Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.48	1.42 ug/l	139049	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	2-Hexanone, 4-hydroxy-3-propyl-	158	C9H18O2	062338-17-4	64
3	Methyl Isobutyl Ketone	100	C6H12O	000108-10-1	58
4	Oxirane, 2-methyl-2-(1-methylethyl)	100	C6H12O	072221-03-5	50



Library Search Compound Report

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Acq On : 10 Oct 2001 3:59 am
Sample : gc/ms unknown
Misc :
MS Integration Params: LSCINT.P

Vial: 16
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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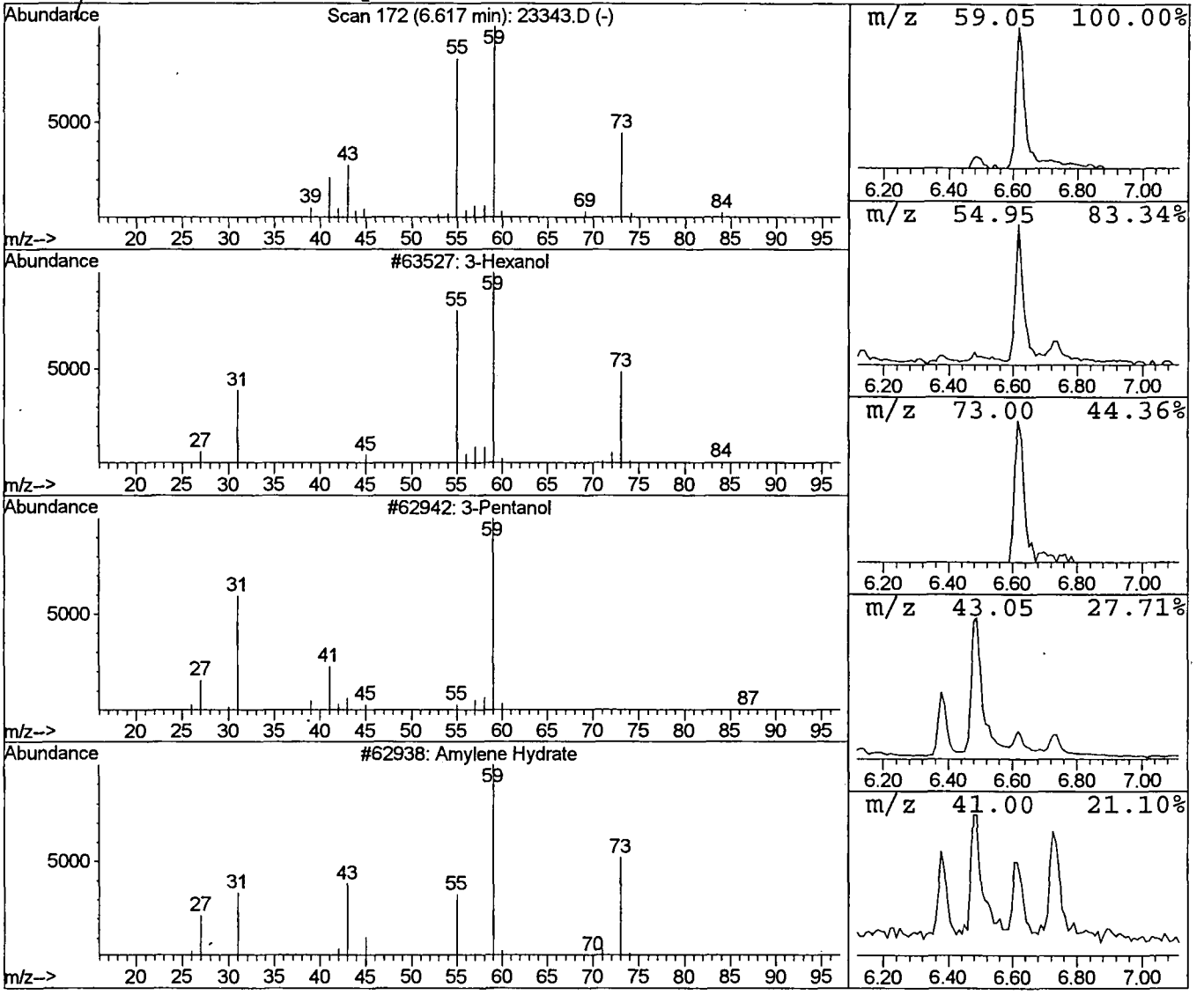
****

Peak Number 3 3-Hexanol Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.62	0.65 ug/l	63357	1,4-Dichlorobenzene-d4	11.76

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	3-Hexanol	102	C6H14O	000623-37-0	83
2	3-Pentanol	88	C5H12O	000584-02-1	42
3	Amylene Hydrate	88	C5H12O	000075-85-4	36
4	3-Pentanol, 2-methyl-	102	C6H14O	000565-67-3	36



Library Search Compound Report

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Acq On : 10 Oct 2001 3:59 am
Sample : gc/ms unknown
Misc :
MS Integration Params: LSCINT.P

Vial: 16
Operator: 209 GALOS
Inst : GC/MS 209
Multiplr: 1.00

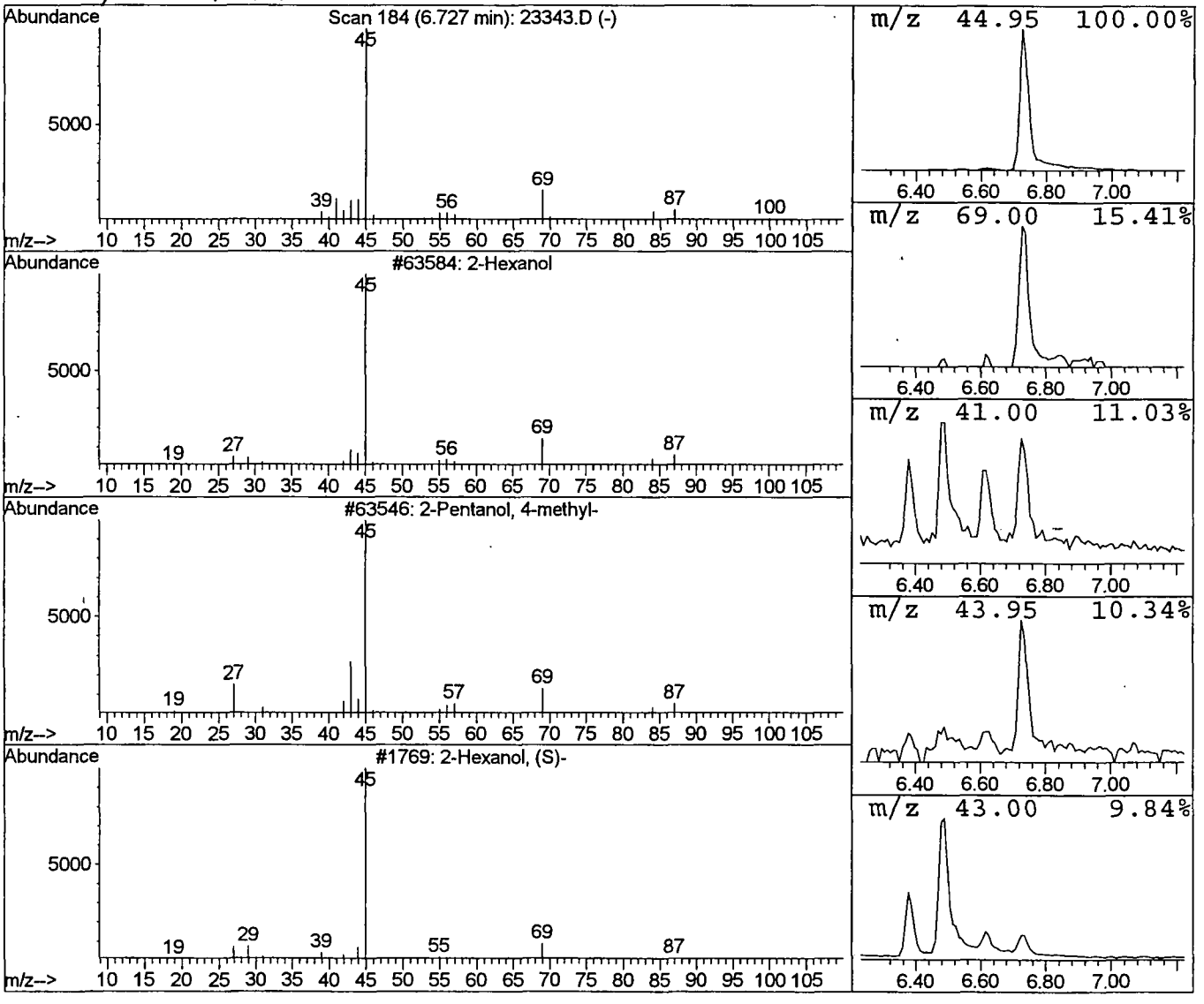
Quant Method : C:\HPCHEM\1\METHODS\NP091101.M (RTE Integrator)
Title : 209 625/TCLP calibration table 7/16/01
Library : C:\DATABASE\NBS75K.L

****

Peak Number 4 2-Hexanol — Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.73	0.92 ug/l	89838	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-Hexanol, (R)-	102	C6H14O	026549-24-6	64



Library Search Compound Report

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Acq On : 10 Oct 2001 3:59 am
Sample : gc/ms unknown
Misc :
MS Integration Params: LSCINT.P

Vial: 16
Operator: 209 GALOS
Inst : GC/MS 209
Multiplr: 1.00

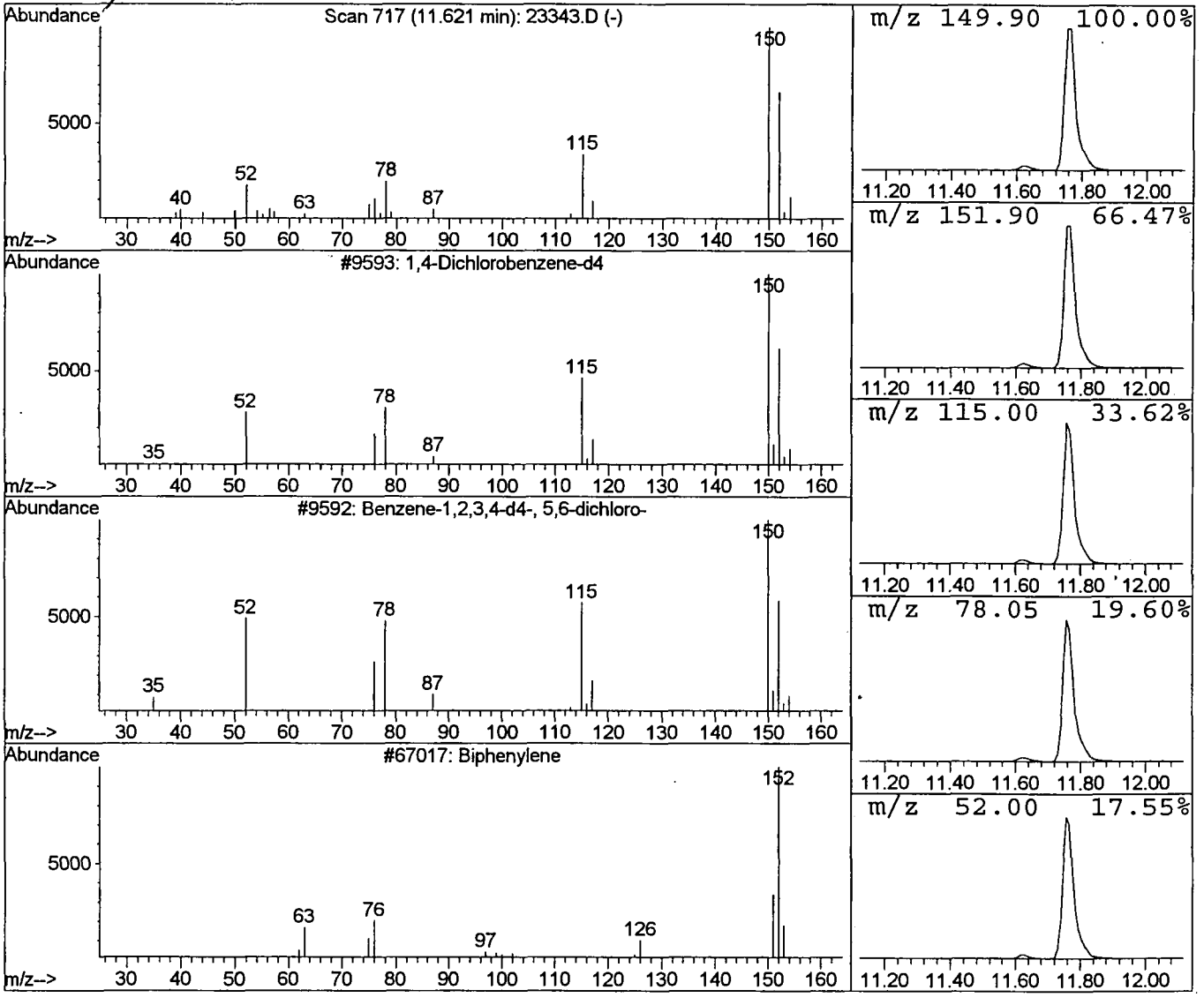
Quant Method : C:\HPCHEM\1\METHODS\NP091101.M (RTE Integrator)
Title : 209 625/TCLP calibration table 7/16/01
Library : C:\DATABASE\NBS75K.L

****

Peak Number 5 1,4-Dichlorobenzene-d4 Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.62	0.54 ug/l	53447	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	64
2	Benzene-1,2,3,4-d4-, 5,6-dichloro-	150	C6Cl2D4	002199-69-1	43
3	Biphenylene	152	C12H8	000259-79-0	9
4	Benzofuran, 7-chloro-	152	C8H5ClO	024410-55-7	9



Library Search Compound Report

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Acq On : 10 Oct 2001 3:59 am
Sample : gc/ms unknown
Misc :
MS Integration Params: LSCINT.P

Vial: 16
Operator: 209 GALOS
Inst : GC/MS 209
Multiplr: 1.00

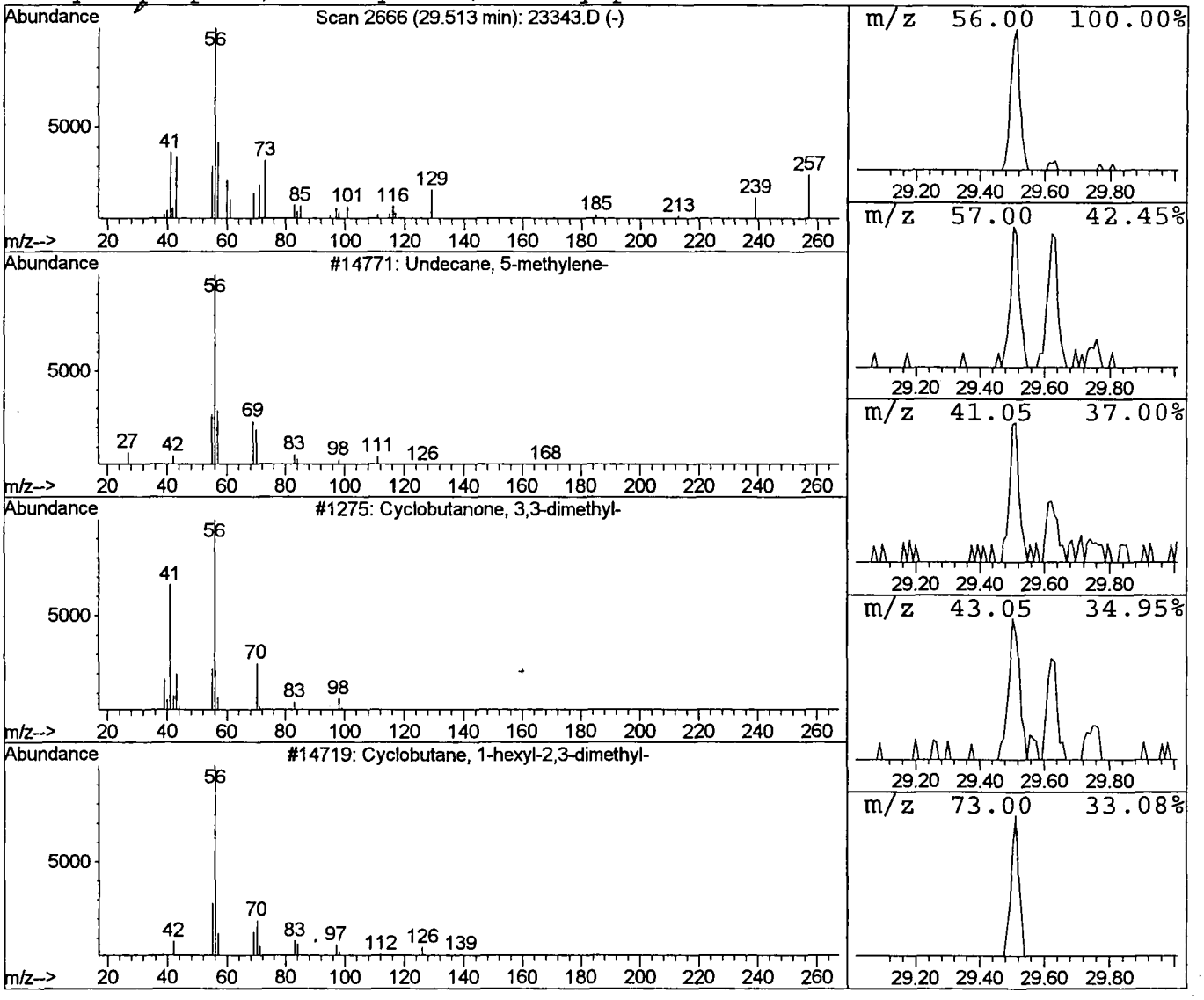
Quant Method : C:\HPCHEM\1\METHODS\NP091101.M (RTE Integrator)
Title : 209 625/TCLP calibration table 7/16/01
Library : C:\DATABASE\NBS75K.L

****

Peak Number 6 Undecane, 5-methylene- Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
29.51	0.41 ug/l	38344	Chrysene-d12	33.12

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Undecane, 5-methylene-	168	C12H24	005698-48-6	27
2	Cyclobutanone, 3,3-dimethyl-	98	C6H10O	001192-33-2	27
3	Cyclobutane, 1-hexyl-2,3-dimethyl-	168	C12H24	055170-84-8	25
4	Cyclopropane, 1-butyl-2-(2-methylpr	154	C11H22	041977-35-9	25



Library Search Compound Report

Data File : C:\HPCHEM\1\DATA\OCT0901\23343.D
Acq On : 10 Oct 2001 3:59 am
Sample : gc/ms unknown
Misc :
MS Integration Params: LSCINT.P

Vial: 16
Operator: 209 GALOS
Inst : GC/MS 209
Multiplr: 1.00

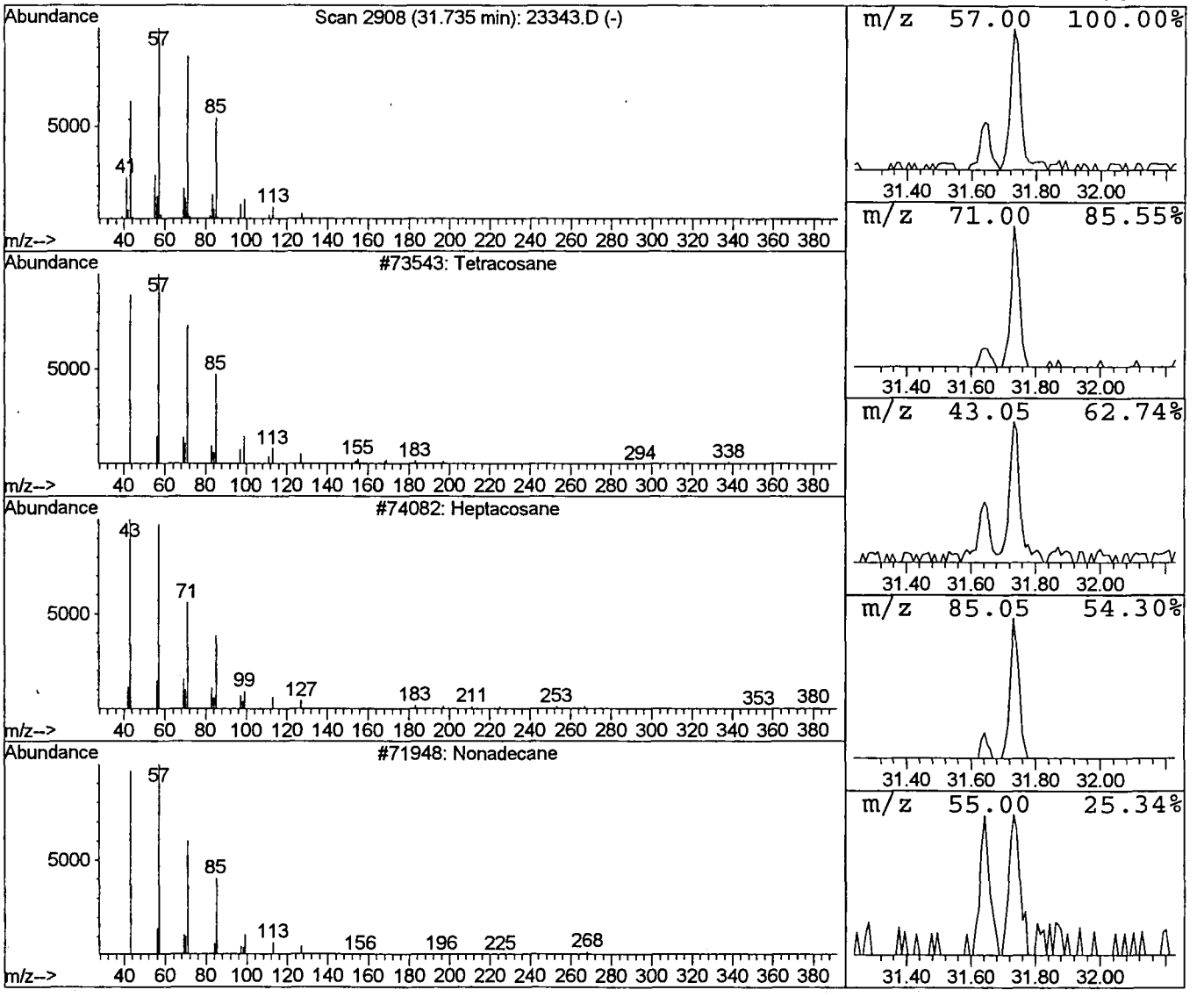
Quant Method : C:\HPCHEM\1\METHODS\NP091101.M (RTE Integrator)
Title : 209 625/TCLP calibration table 7/16/01
Library : C:\DATABASE\NBS75K.L

****

Peak Number 7 Tetracosane — Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
31.73	0.45 ug/l	41539	Chrysene-d12	33.12

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Tetracosane	338	C24H50	000646-31-1	90
2	Heptacosane	380	C27H56	000593-49-7	86
3	Nonadecane	268	C19H40	000629-92-5	78
4	Pentacosane	352	C25H52	000629-99-2	78



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Acq On : 10 Oct 2001 3:59 am
Sample : gc/ms unknown
Misc :
MS Integration Params: LSCINT.P

Vial: 16
Operator: 209 GALOS
Inst : GC/MS 209
Multiplr: 1.00

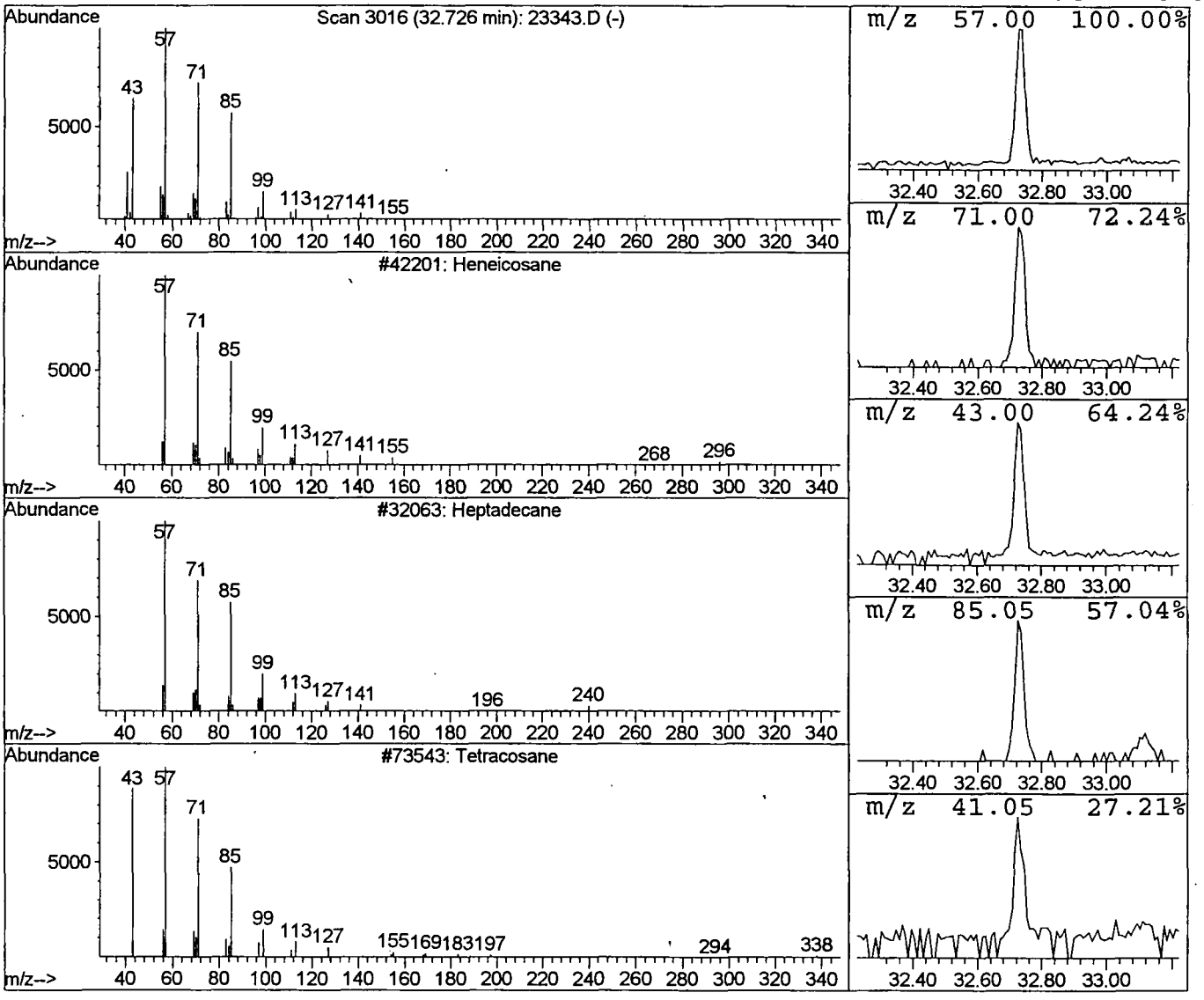
Quant Method : C:\HPCHEM\1\METHODS\NP091101.M (RTE Integrator)
Title : 209 625/TCLP calibration table 7/16/01
Library : C:\DATABASE\NBS75K.L

****

Peak Number 8 Heneicosane Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
32.73	0.53 ug/l	49818	Chrysene-d12	33.12

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heneicosane	296	C21H44	000629-94-7	91
2	Heptadecane	240	C17H36	000629-78-7	90
3	Tetracosane	338	C24H50	000646-31-1	87
4	Octadecane	254	C18H38	000593-45-3	86



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MS Integration Params: LSCINT.P

Vial: 16
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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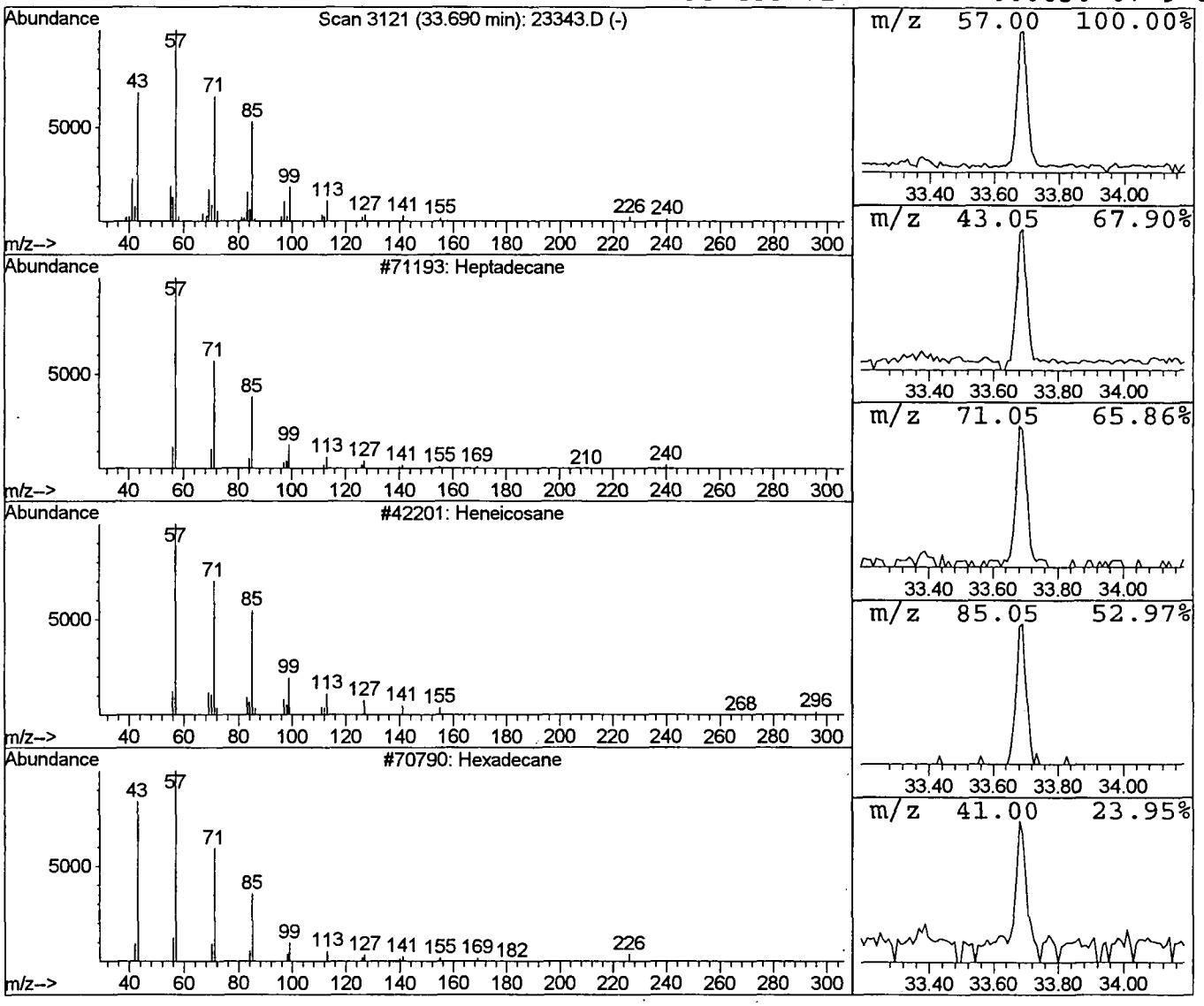
Peak Number 9 Heptadecane — Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
33.69	0.73 ug/l	67583	Chrysene-d12	33.12

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heptadecane	240	C17H36	000629-78-7	94
2	Heneicosane	296	C21H44	000629-94-7	91
3	Hexadecane	226	C16H34	000544-76-3	90
4	Pentatriacontane	493	C35H72	000630-07-9	83



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Vial: 16
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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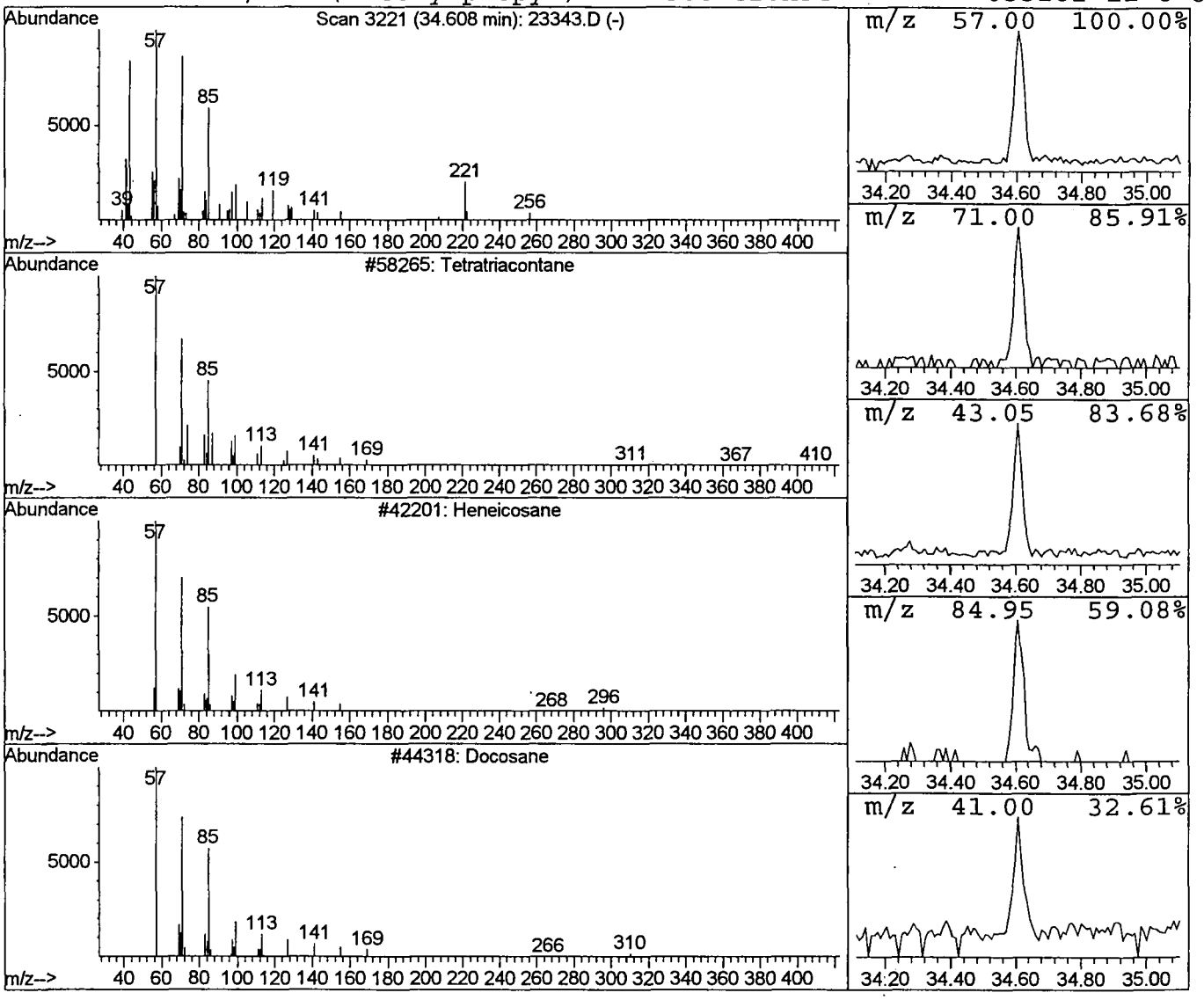
Peak Number 10 Tetratriacontane Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
34.61	0.53 ug/l	49254	Chrysene-d12	33.12

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Tetratriacontane	479	C34H70	014167-59-0	83
2	Heneicosane	296	C21H44	000629-94-7	80
3	Docosane	310	C22H46	000629-97-0	80
4	Heneicosane, 11-(1-ethylpropyl)-	366	C26H54	055282-11-6	80



Library Search Compound Report

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Vial: 16
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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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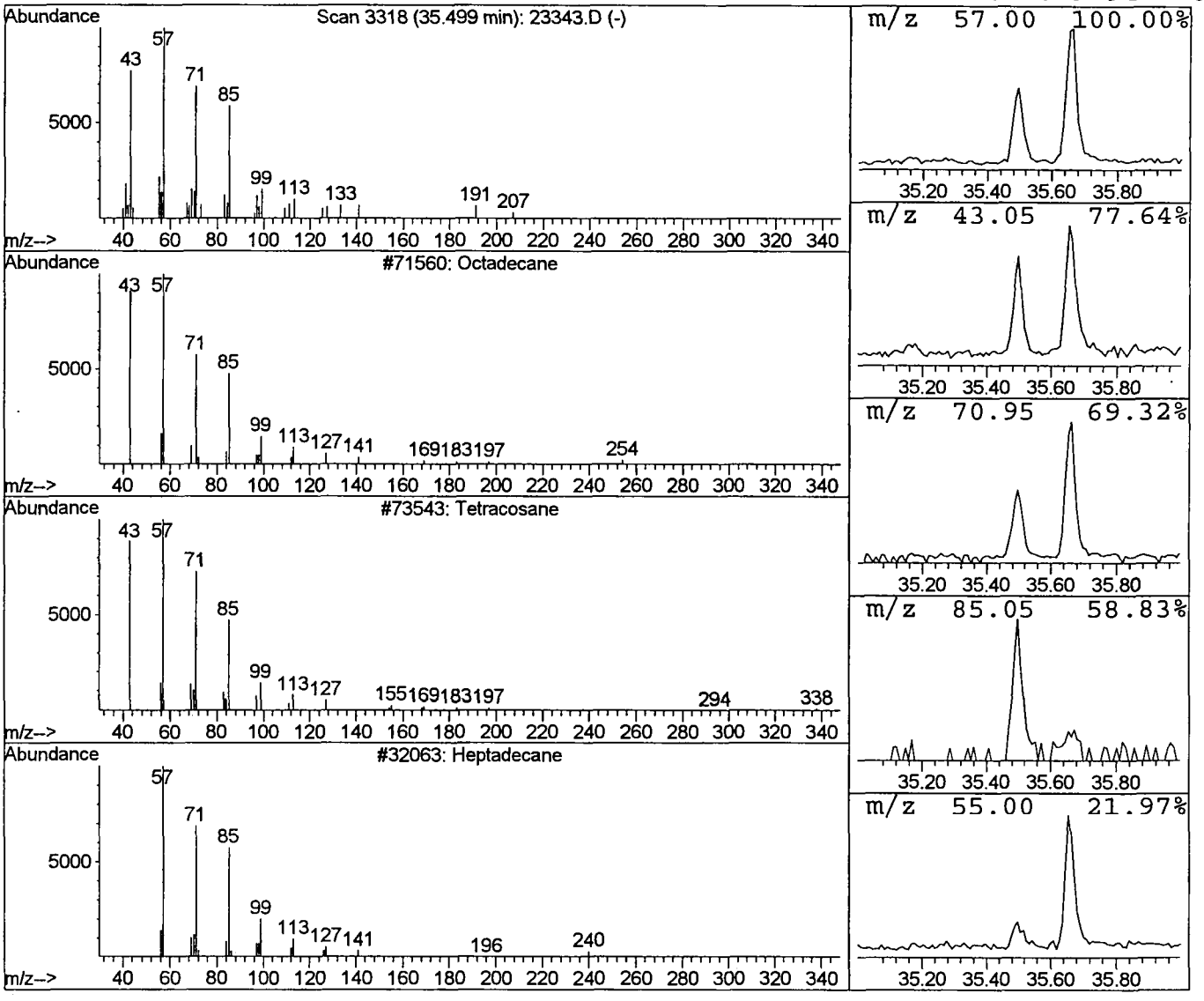
Peak Number 11 Octadecane — Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
35.50	0.43 ug/l	35735	Perylene-d12	37.23

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Octadecane	254	C18H38	000593-45-3	86
2	Tetracosane	338	C24H50	000646-31-1	86
3	Heptadecane	240	C17H36	000629-78-7	86
4	Heneicosane	296	C21H44	000629-94-7	86



Library Search Compound Report

Data File : C:\HPCHEM\1\DATA\OCT0901\23343.D
Acq On : 10 Oct 2001 3:59 am
Sample : gc/ms unknown
Misc :
MS Integration Params: LSCINT.P

Vial: 16
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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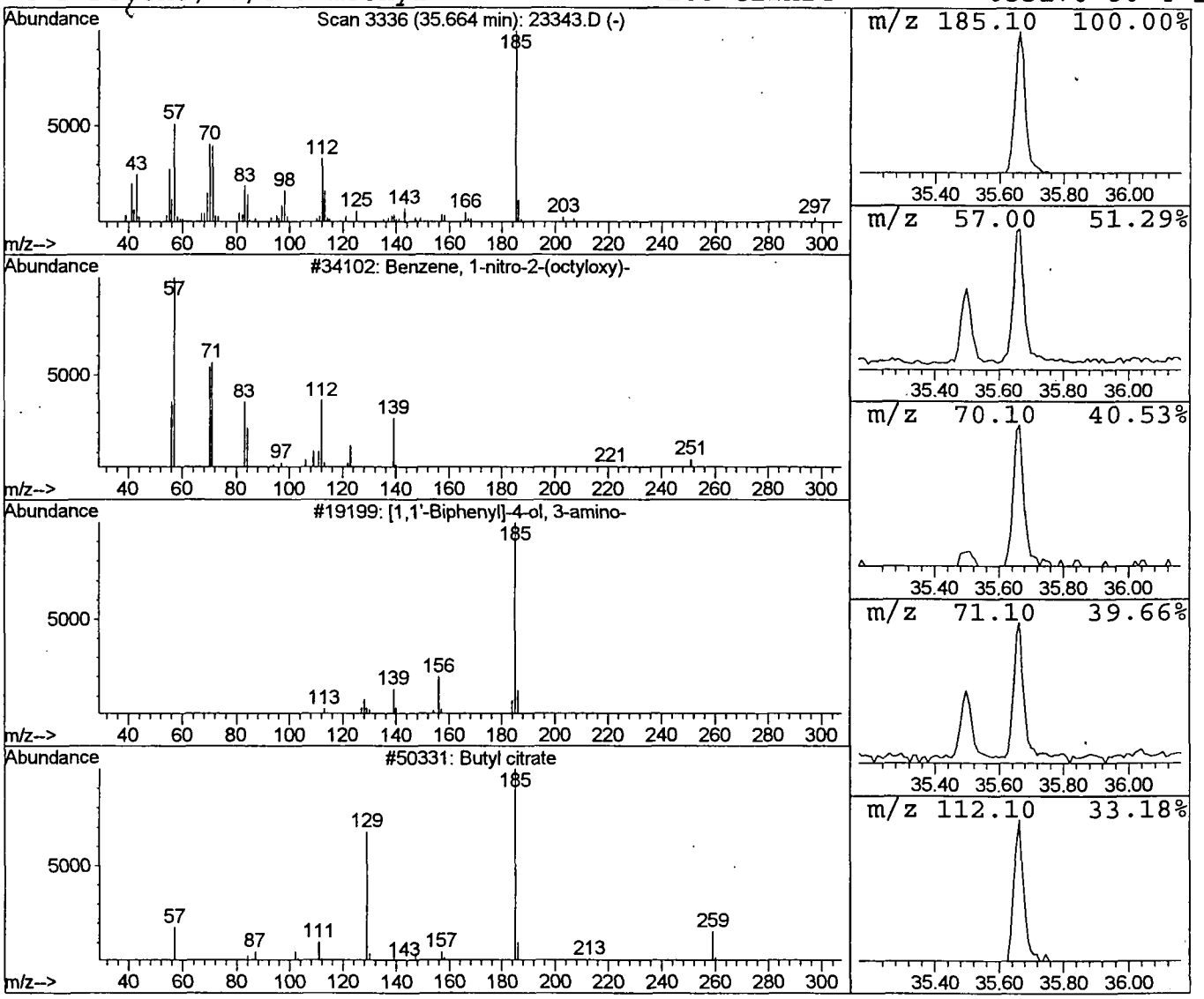
Peak Number 12 Benzene, 1-nitro-2-(octyloxy)- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
35.66	1.64 ug/l	135080	Perylene-d12	37.23

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1-nitro-2-(octyloxy)-	251	C14H21NO3	037682-29-4	17
2	[1,1'-Biphenyl]-4-ol, 3-amino-	185	C12H11NO	001134-36-7	16
3	Butyl citrate	360	C18H32O7	000077-94-1	16
4	1-Decene, 2,4-dimethyl-	168	C12H24	055170-80-4	16



Tentatively Identified Compound (LSC) summary

Operator ID: 209 GALOS Date Acquired: 10 Oct 2001 3:59 am
Data File: C:\HPCHEM\1\DATA\OCT0901\23343.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	0.6	ug/l	63697	ISTD01	11.76	1962260	20.0
2-Hexanone	6.48	1.4	ug/l	139049	ISTD01	11.76	1962260	20.0
3-Hexanol	6.62	0.6	ug/l	63357	ISTD01	11.76	1962260	20.0
2-Hexanol	6.73	0.9	ug/l	89838	ISTD01	11.76	1962260	20.0
1,4-Dichlorobenzene-	11.62	0.5	ug/l	53447	ISTD01	11.76	1962260	20.0
Undecane, 5-methylen	29.51	0.4	ug/l	38344	ISTD05	33.12	1862630	20.0
Tetracosane	31.73	0.4	ug/l	41539	ISTD05	33.12	1862630	20.0
Heneicosane	32.73	0.5	ug/l	49818	ISTD05	33.12	1862630	20.0
Heptadecane	33.69	0.7	ug/l	67583	ISTD05	33.12	1862630	20.0
Tetratriacontane	34.61	0.5	ug/l	49254	ISTD05	33.12	1862630	20.0
Octadecane	35.50	0.4	ug/l	35735	ISTD06	37.23	1645630	20.0
Benzene, 1-nitro-2-(	35.66	1.6	ug/l	135080	ISTD06	37.23	1645630	20.0

23343.D NP091101.M Thu Oct 11 13:05:51 2001

Comments for Sample AD23343 - Analysis $GCMS

Westchester Dept. of Labs & Research
User: Vinci, David L
Date: 10-19-2001 Time: 15:34:19

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. An investigation into the source of these compounds in the Laboratory Blanks and also in many of these NYCDEP samples is still under way. An advisory will be made as soon as the investigation is concluded.