User: Galos, Marie
Westchester Dept. of Labs & Research
Date: 11/01/2001 Time: 13:52:24

Sample: AD23839
Analysis: $GCMS GCMS Scan For Unknowns
Units: ug
Started: 11/1/01 13:52 Ended: 11/1/01 13:52 Analyst: MG
Page: 1

	Component Name	Vial	Result
1	Other unknown compounds		see comment

NYC 9/11 Public Portal Document

Quantitation Report (QT Reviewed)

Data File : C:\HPCHEM\1\DATA\OCT3001\23839.D Vial: 13
Acq On : 31 Oct 2001 2:14 am Operator: 209 GALOS
Sample : gc/ms unknown Inst : GC/MS 209
Misc : Multiplr: 1.00
MS Integration Params: RTEINT.P
Quant Time: Oct 31 12:17 2001 Quant Results File: NP101901.RES

Quant Method : C:\HPCHEM\1\METHODS\NP101901.M (RTE Integrator)
Title : 209 625/TCLP calibration table
Last Update : Tue Oct 23 10:12:13 2001
Response via : Initial Calibration
DataAcq Meth : NP101901

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	11.72	152	988170	20.00	ug/l	-0.04
19) Naphthalene-d8	15.32	136	3866423	20.00	ug/l	-0.04
31) Acenaphthene-d10	20.60	164	1822500	20.00	ug/l	-0.03
49) Phenanthrene-d10	25.03	188	2556532	20.00	ug/l	-0.03
59) Chrysene-d12	33.06	240	1503579	20.00	ug/l	-0.04
68) Perylene-d12	37.16	264	1044160	20.00	ug/l	-0.03
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	11.72	132	891	0.01	ug/l	0.45
Spiked Amount 75.000			Recovery =	0.01%		
7) 1,2-Dichlorobenzene-d4	12.23	152	11231	0.16	ug/l	-0.04
Spiked Amount 50.000			Recovery =	0.32%		
20) Nitrobenzene-d5	0.00	82	0	0.00	ug/l	
Spiked Amount 50.000			Recovery =	0.00%		
32) 2-Fluorobiphenyl	18.74	172	3922	0.02	ug/l	0.09
Spiked Amount 50.000			Recovery =	0.04%		
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

NYC-WTC_000089367

NYC 9/11 Public Portal Document

Quantitation Report (QT Reviewed)

Data File : C:\HPCHEM\1\DATA\OCT3001\23839.D Vial: 13
Acq On : 31 Oct 2001 2:14 am Operator: 209 GALOS
Sample : gc/ms unknown Inst : GC/MS 209
Misc : Multiplr: 1.00
MS Integration Params: RTEINT.P
Quant Time: Oct 31 12:17 2001 Quant Results File: NP101901.RES

Quant Method : C:\HPCHEM\1\METHODS\NP101901.M (RTE Integrator)
Title : 209 625/TCLP calibration table
Last Update : Tue Oct 23 10:12:13 2001
Response via : Initial Calibration
DataAcq Meth : NP101901

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	15.37	128	1091	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	21.11	165	198	N.D.	
45) Diethylphthalate	22.11	149	457	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	25.08	178	1315	N.D.	
56) Anthracene	25.08	178	1315	N.D.	
57) Di-n-butylphthalate	27.03	149	5791	N.D.	
58) Fluoranthene	28.71	202	330	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	33.06	228	4460	N.D.	
66) Bis(2-ethylhexyl)phthalate	33.34	149	2969	N.D.	
67) Chrysene	33.06	228	4460	N.D.	
69) Di-n-Octylphthalate	35.09	149	534	N.D.	
70) Benzo(b)fluoranthene	36.14	252	2281	N.D.	

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

NYC-WTC_000089368

NYC 9/11 Public Portal Document

Quantitation Report (QT Reviewed)

Data File : C:\HPCHEM\1\DATA\OCT3001\23839.D Vial: 13
Acq On : 31 Oct 2001 2:14 am Operator: 209 GALOS
Sample : gc/ms unknown Inst : GC/MS 209
Misc : Multiplr: 1.00
MS Integration Params: RTEINT.P
Quant Time: Oct 31 12:17 2001 Quant Results File: NP101901.RES

Quant Method : C:\HPCHEM\1\METHODS\NP101901.M (RTE Integrator)
Title : 209 625/TCLP calibration table
Last Update : Tue Oct 23 10:12:13 2001
Response via : Initial Calibration
DataAcq Meth : NP101901

Compound	R.T.	QIon	Response	Conc Unit	Qvalue
71) Benzo(k)fluoranthene	36.14	252	1439	N.D.	
72) Benzo(a)pyrene	36.98	252	825	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
23839.D NP101901.M Wed Oct 31 12:18:04 2001 Page 3

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NYC 9/11 Public Portal Document

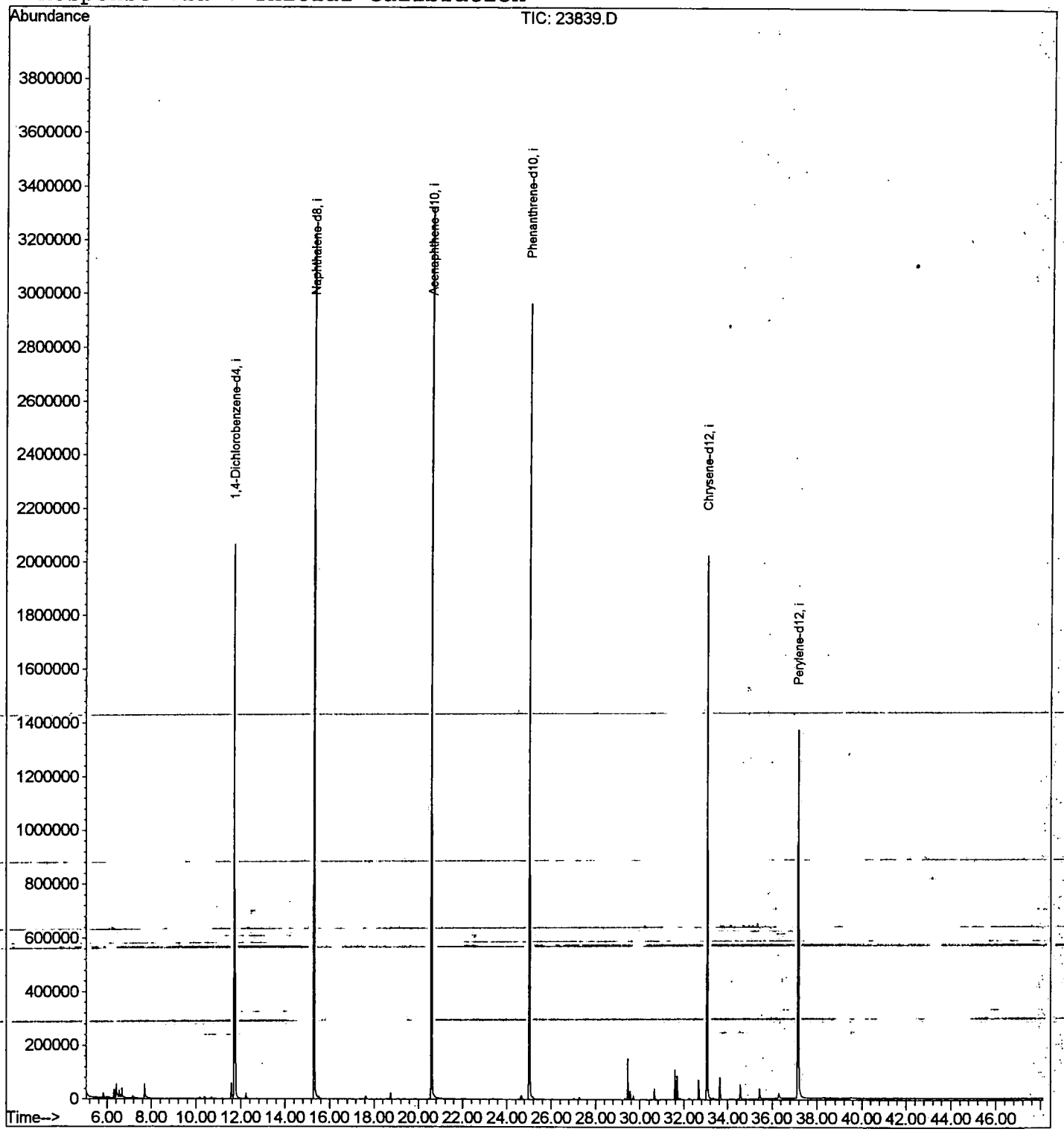
Quantitation Report

Data File : C:\HPCHEM\1\DATA\OCT3001\23839.D
Acq On : 31 Oct 2001 2:14 am
Sample : gc/ms unknown
Misc :
MS Integration Params: RTEINT.P
Quant Time: Oct 31 12:17 2001

Vial: 13
Operator: 209 GALOS
Inst : GC/MS 209
Multiplr: 1.00

Quant Results File: NP101901.RES

Method : C:\HPCHEM\1\METHODS\NP101901.M (RTE Integrator)
Title : 209 625/TCLP calibration table
Last Update : Tue Oct 23 10:12:13 2001
Response via : Initial Calibration



NYC-WTC_000089370

NYC 9/11 Public Portal Document

LSC Area Percent Report

Data File : C:\HPCHEM\1\DATA\OCT3001\23839.D
Acq On : 31 Oct 2001 2:14 am
Sample : gc/ms unknown
Misc :
MS Integration Params: LSCINT.P

Vial: 13
Operator: 209 GALOS
Inst : GC/MS 209
Multiplr: 1.00

Method : C:\HPCHEM\1\METHODS\NP103001.M (RTE Integrator)
Title : 209 625/TCLP calibration table
Smoothing : OFF
Sampling : 1
Start Thrs: 0.2
Stop Thrs : 0

Filtering: 5
Min Area: 1 % of largest Peak
Max Peaks: 100
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	6.428	147	151	161	rVV2	51307	124456	1.72%	0.343%
2	7.695	286	289	302	rBV	53677	144202	2.00%	0.397%
3	11.569	705	711	721	rBV	59323	144313	2.00%	0.398%
4	11.716	721	727	746	rVV	2064616	5165664	71.55%	14.229%
5	15.324	1113	1120	1137	rBV	3225392	7037599	97.47%	19.385%
6	20.602	1688	1695	1712	rBV	3326722	7219922	100.00%	19.887%
7	25.027	2170	2177	2189	rBV	2967093	6370160	88.23%	17.546%
8	29.452	2654	2659	2667	rBV	152359	291219	4.03%	0.802%
9	30.646	2784	2789	2795	rBV	41398	82153	1.14%	0.226%
10	31.591	2884	2892	2898	rBV	111157	224590	3.11%	0.619%
11	31.683	2898	2902	2912	rVB	87817	178368	2.47%	0.491%
12	32.674	3003	3010	3019	rBV	73383	164261	2.28%	0.452%
13	33.060	3044	3052	3066	rBV2	2029873	4798893	66.47%	13.218%
14	33.638	3110	3115	3124	rBV	80929	185301	2.57%	0.510%
15	34.556	3210	3215	3228	rVB	56463	129695	1.80%	0.357%
16	35.447	3306	3312	3325	rBV2	39882	110329	1.53%	0.304%
17	37.164	3490	3499	3523	rBV2	1372668	3933742	54.48%	10.835%

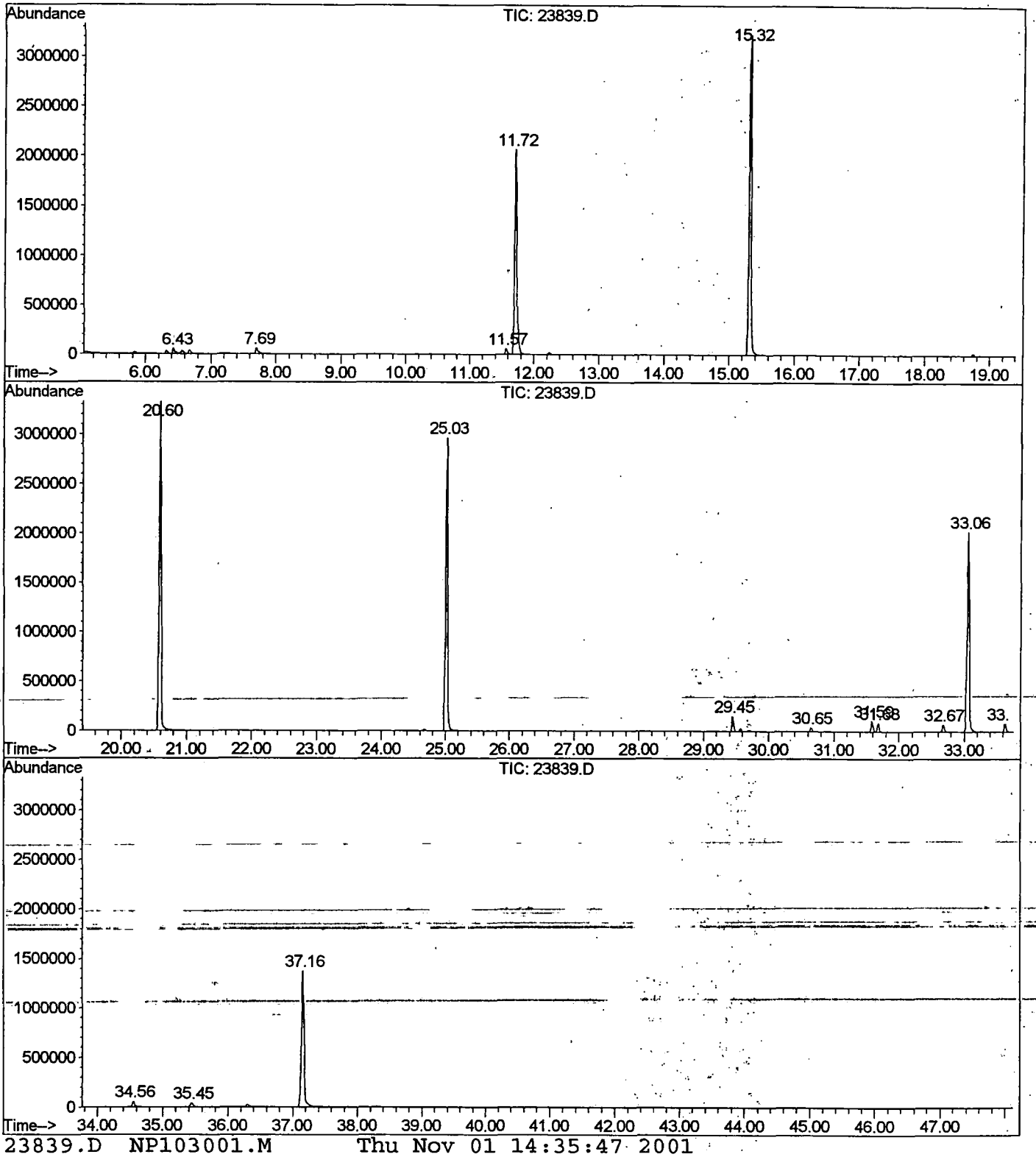
Sum of corrected areas: 36304867

23839.D NP103001.M Thu Nov 01 14:35:44 2001

NYC 9/11 Public Portal Document

LSC Report - Integrated Chromatogram

File : C:\HPCHEM\1\DATA\OCT3001\23839.D
Operator : 209 GALOS
Acquired : 31 Oct 2001 2:14 am using AcqMethod NP101901
Instrument : GC/MS 209
Sample Name: gc/ms unknown
Misc Info :
Vial Number: 13
Quant File :NP101901.RES (RTE Integrator)



23839.D NP103001.M Thu Nov 01 14:35:47 2001

NYC-WTC_000089372

NYC 9/11 Public Portal Document

Library Search Compound Report

Data File : C:\HPCHEM\1\DATA\OCT3001\23839.D
Acq On : 31 Oct 2001 2:14 am
Sample : gc/ms unknown
Misc :
MS Integration Params: LSCINT.P

Vial: 13
Operator: 209 GALOS
Inst : GC/MS 209
Multiplr: 1.00

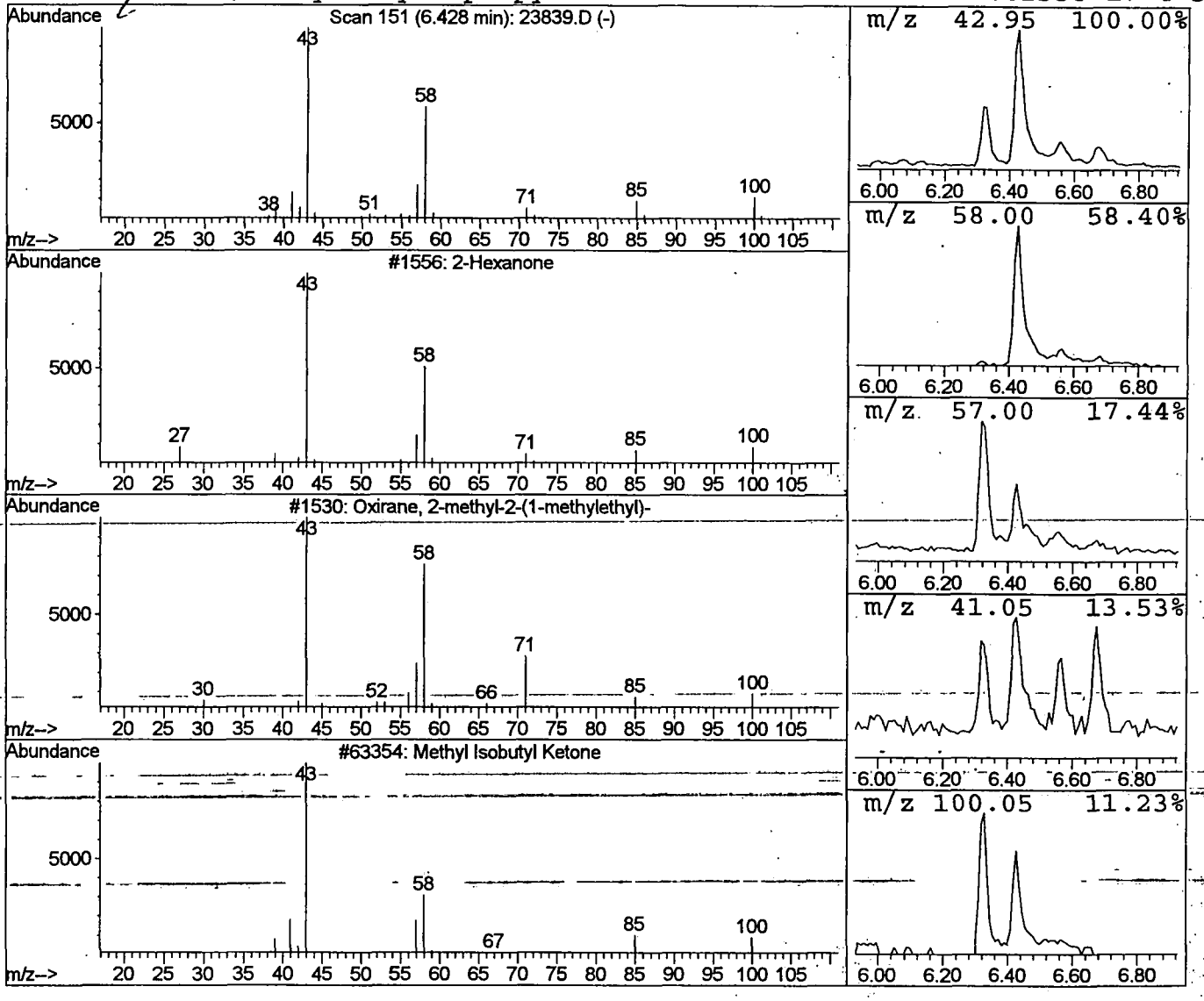
Quant Method : C:\HPCHEM\1\METHODS\NP101901.M (RTE Integrator)
Title : 209 625/TCLP calibration table
Library : C:\DATABASE\NBS75K.L

****

Peak Number 1 2-Hexanone Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.43	0.48 ug/l	124456	1,4-Dichlorobenzene-d4	11.72

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	80
3	Methyl Isobutyl Ketone	100	C6H12O	000108-10-1	58
4	2-Hexanone, 4-hydroxy-3-propyl-	158	C9H18O2	062338-17-4	56



NYC-WTC_000089373

NYC 9/11 Public Portal Document

Library Search Compound Report

Data File : C:\HPCHEM\1\DATA\OCT3001\23839.D
Acq On : 31 Oct 2001 2:14 am
Sample : gc/ms unknown
Misc :
MS Integration Params: LSCINT.P

Vial: 13
Operator: 209 GALOS
Inst : GC/MS 209
Multiplr: 1.00

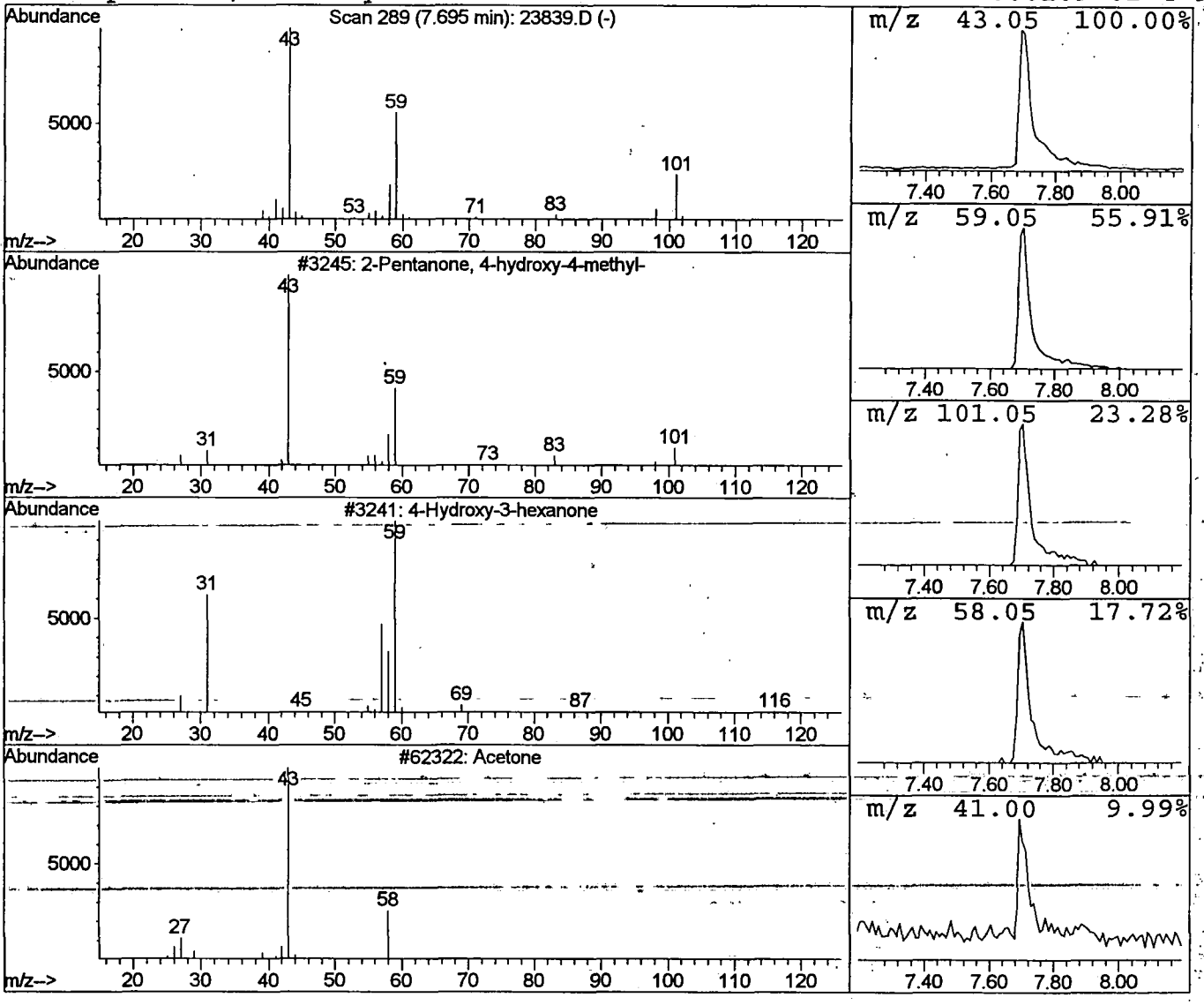
Quant Method : C:\HPCHEM\1\METHODS\NP101901.M (RTE Integrator)
Title : 209 625/TCLP calibration table
Library : C:\DATABASE\NBS75K.L

****

Peak Number 2 2-Pentanone, 4-hydroxy-4-methy Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.69	0.56 ug/l	144202	1,4-Dichlorobenzene-d4	11.72

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	50
2	4-Hydroxy-3-hexanone	116	C6H12O2	004984-85-4	17
3	Acetone	58	C3H6O	000067-64-1	12
4	Morpholine, 4-methyl-	101	C5H11NO	000109-02-4	9



NYC-WTC_000089374

Library Search Compound Report

Data File : C:\HPCHEM\1\DATA\OCT3001\23839.D
Acq On : 31 Oct 2001 2:14 am
Sample : gc/ms unknown
Misc :
MS Integration Params: LSCINT.P

Vial: 13
Operator: 209 GALOS
Inst : GC/MS 209
Multiplr: 1.00

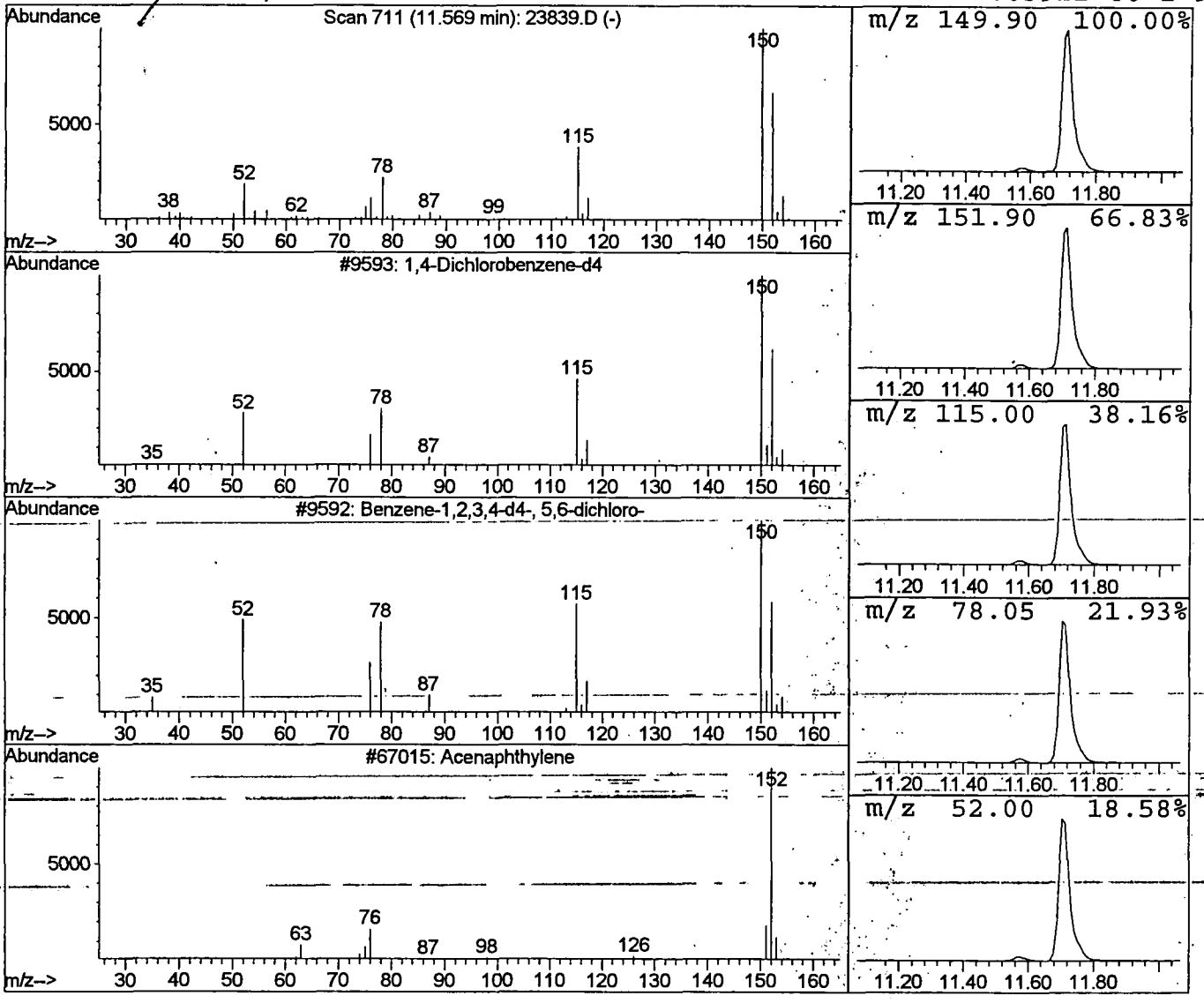
Quant Method : C:\HPCHEM\1\METHODS\NP101901.M (RTE Integrator)
Title : 209 625/TCLP calibration table
Library : C:\DATABASE\NBS75K.L

****

Peak Number 3 1,4-Dichlorobenzene-d4 Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.57	0.56 ug/l	144313	1,4-Dichlorobenzene-d4	11.72

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,4-Dichlorobenzene-d4	150	C6Cl2D4	003855-82-1	94
2	Benzene-1,2,3,4-d4-, 5,6-dichloro-	150	C6Cl2D4	002199-69-1	38
3	Acenaphthylene	152	C12H8	000208-96-8	10
4	Benzonitrile, 2-amino-5-chloro-	152	C7H5ClN2	005922-60-1	9



NYC 9/11 Public Portal Document

Library Search Compound Report

Data File : C:\HPCHEM\1\DATA\OCT3001\23839.D
Acq On : 31 Oct 2001 2:14 am
Sample : gc/ms unknown
Misc :
MS Integration Params: LSCINT.P

Vial: 13
Operator: 209 GALOS
Inst : GC/MS 209
Multiplr: 1.00

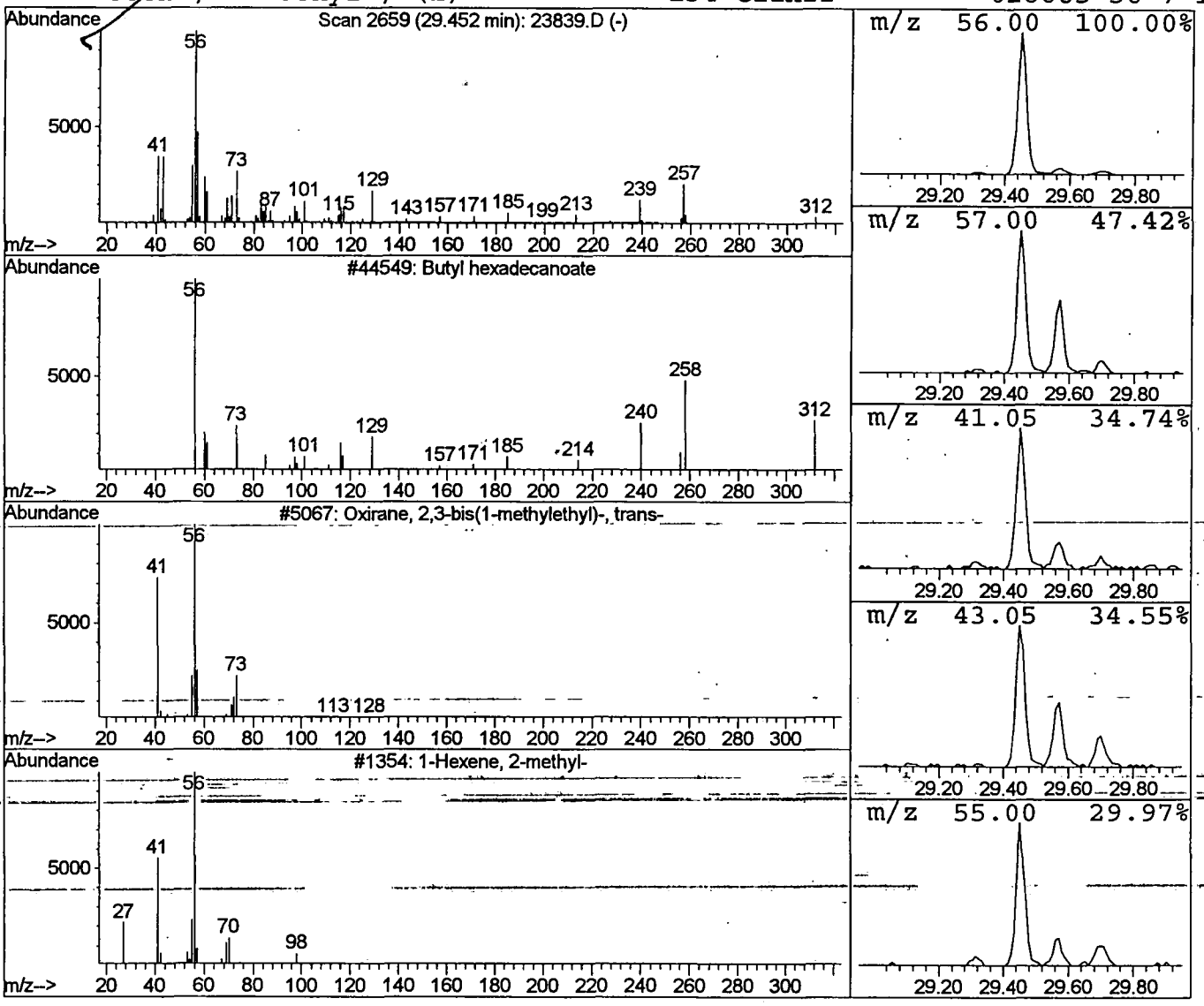
Quant Method : C:\HPCHEM\1\METHODS\NP101901.M (RTE Integrator)
Title : 209 625/TCLP calibration table
Library : C:\DATABASE\NBS75K.L

****

Peak Number 4 Butyl hexadecanoate Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
29.45	1.21 ug/l	291219	Chrysene-d12	33.06

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Butyl hexadecanoate	312	C20H40O2	000000-00-0	35
2	Oxirane, 2,3-bis(1-methylethyl)-, t	128	C8H16O	054644-32-5	32
3	1-Hexene, 2-methyl-	98	C7H14	006094-02-6	27
4	4-Decene, 2-methyl-, (E)-	154	C11H22	028665-56-7	16



NYC-WTC_000089376

NYC 9/11 Public Portal Document

Library Search Compound Report

Data File : C:\HPCHEM\1\DATA\OCT3001\23839.D
Acq On : 31 Oct 2001 2:14 am
Sample : gc/ms unknown
Misc :
MS Integration Params: LSCINT.P

Vial: 13
Operator: 209 GALOS
Inst : GC/MS 209
Multiplr: 1.00

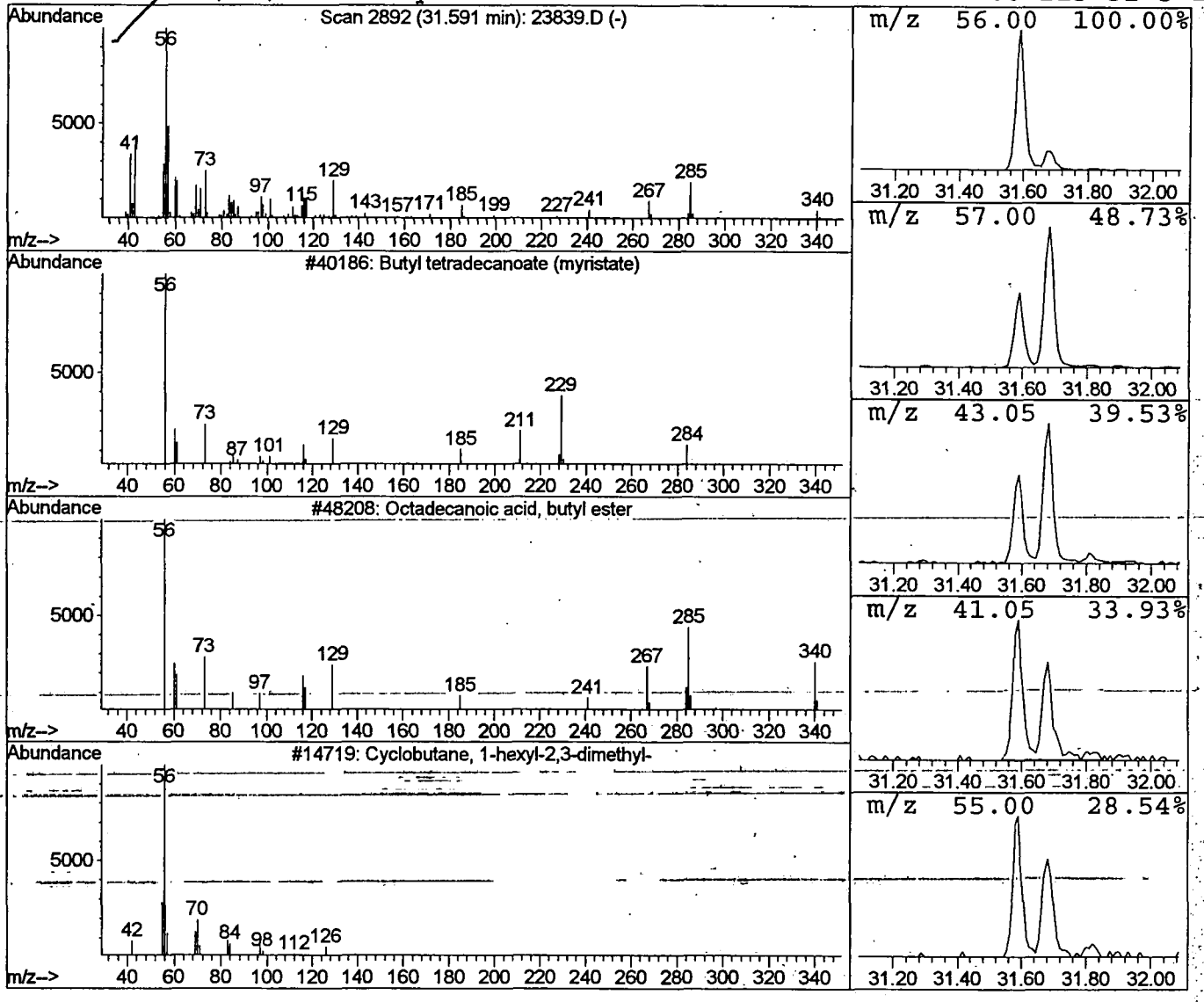
Quant Method : C:\HPCHEM\1\METHODS\NP101901.M (RTE Integrator)
Title : 209 625/TCLP calibration table
Library : C:\DATABASE\NBS75K.L

****

Peak Number 5 Butyl tetradecanoate (myristat Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
31.59	0.94 ug/l	224590	Chrysene-d12	33.06

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Butyl tetradecanoate (myristate)	284	C18H36O2	000000-00-0	72
2	Octadecanoic acid, butyl ester	340	C22H44O2	000123-95-5	50
3	Cyclobutane, 1-hexyl-2,3-dimethyl-	168	C12H24	055170-84-8	27
4	1-Pentene, 2,4-dimethyl-	98	C7H14	002213-32-3	27



NYC-WTC_000089377

Library Search Compound Report

Data File : C:\HPCHEM\1\DATA\OCT3001\23839.D
Acq On : 31 Oct 2001 2:14 am
Sample : gc/ms unknown
Misc :
MS Integration Params: LSCINT.P

Vial: 13
Operator: 209 GALOS
Inst : GC/MS 209
Multiplr: 1.00

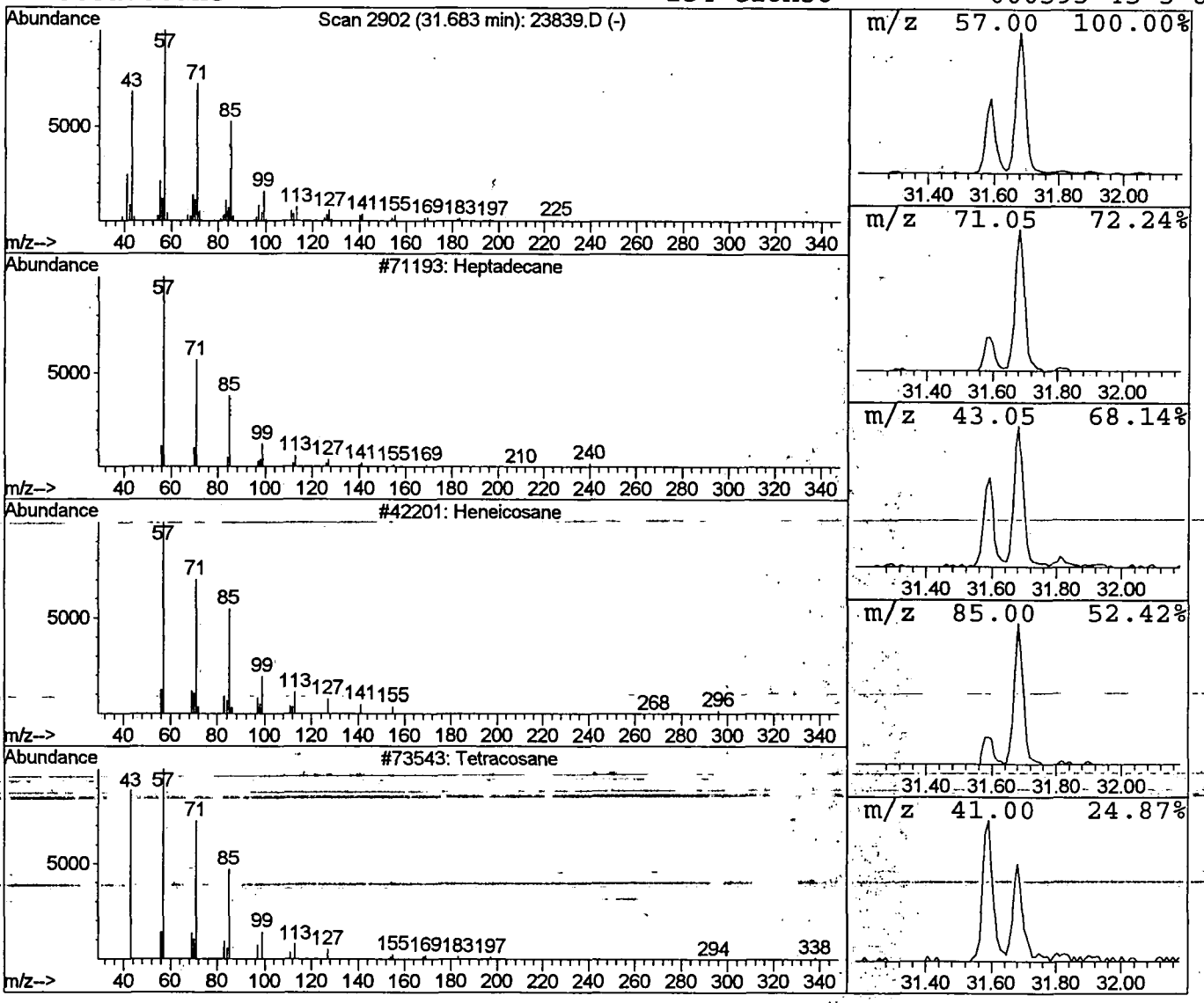
Quant Method : C:\HPCHEM\1\METHODS\NP101901.M (RTE Integrator)
Title : 209 625/TCLP calibration table
Library : C:\DATABASE\NBS75K.L

****

Peak Number 6 Heptadecane — Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
31.68	0.74 ug/l	178368	Chrysene-d12	33.06

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heptadecane	240	C17H36	000629-78-7	91
2	Heneicosane	296	C21H44	000629-94-7	90
3	Tetracosane	338	C24H50	000646-31-1	87
4	Octadecane	254	C18H38	000593-45-3	83



NYC 9/11 Public Portal Document

Library Search Compound Report

Data File : C:\HPCHEM\1\DATA\OCT3001\23839.D
Acq On : 31 Oct 2001 2:14 am
Sample : gc/ms unknown
Misc :
MS Integration Params: LSCINT.P

Vial: 13
Operator: 209 GALOS
Inst : GC/MS 209
Multiplr: 1.00

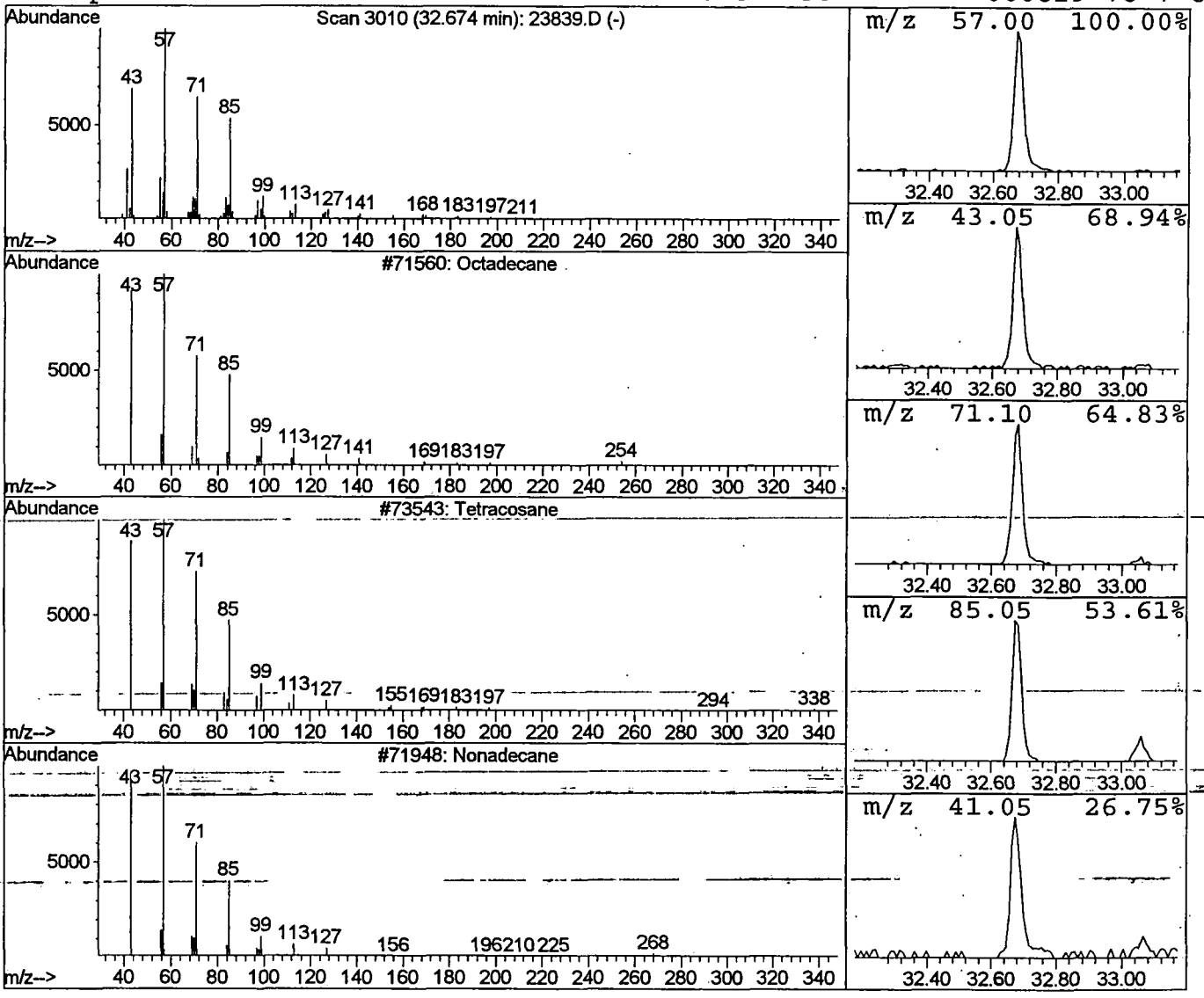
Quant Method : C:\HPCHEM\1\METHODS\NP101901.M (RTE Integrator)
Title : 209 625/TCLP calibration table
Library : C:\DATABASE\NBS75K.L

****

Peak Number 7 Octadecane — Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
32.67	0.68 ug/l	164261	Chrysene-d12	33.06

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Octadecane	254	C18H38	000593-45-3	90
2	Tetracosane	338	C24H50	000646-31-1	90
3	Nonadecane	268	C19H40	000629-92-5	90
4	Heptadecane	240	C17H36	000629-78-7	87



NYC-WTC_000089379

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Library Search Compound Report

Data File : C:\HPCHEM\1\DATA\OCT3001\23839.D
Acq On : 31 Oct 2001 2:14 am
Sample : gc/ms unknown
Misc :
MS Integration Params: LSCINT.P

Vial: 13
Operator: 209 GALOS
Inst : GC/MS 209
Multiplr: 1.00

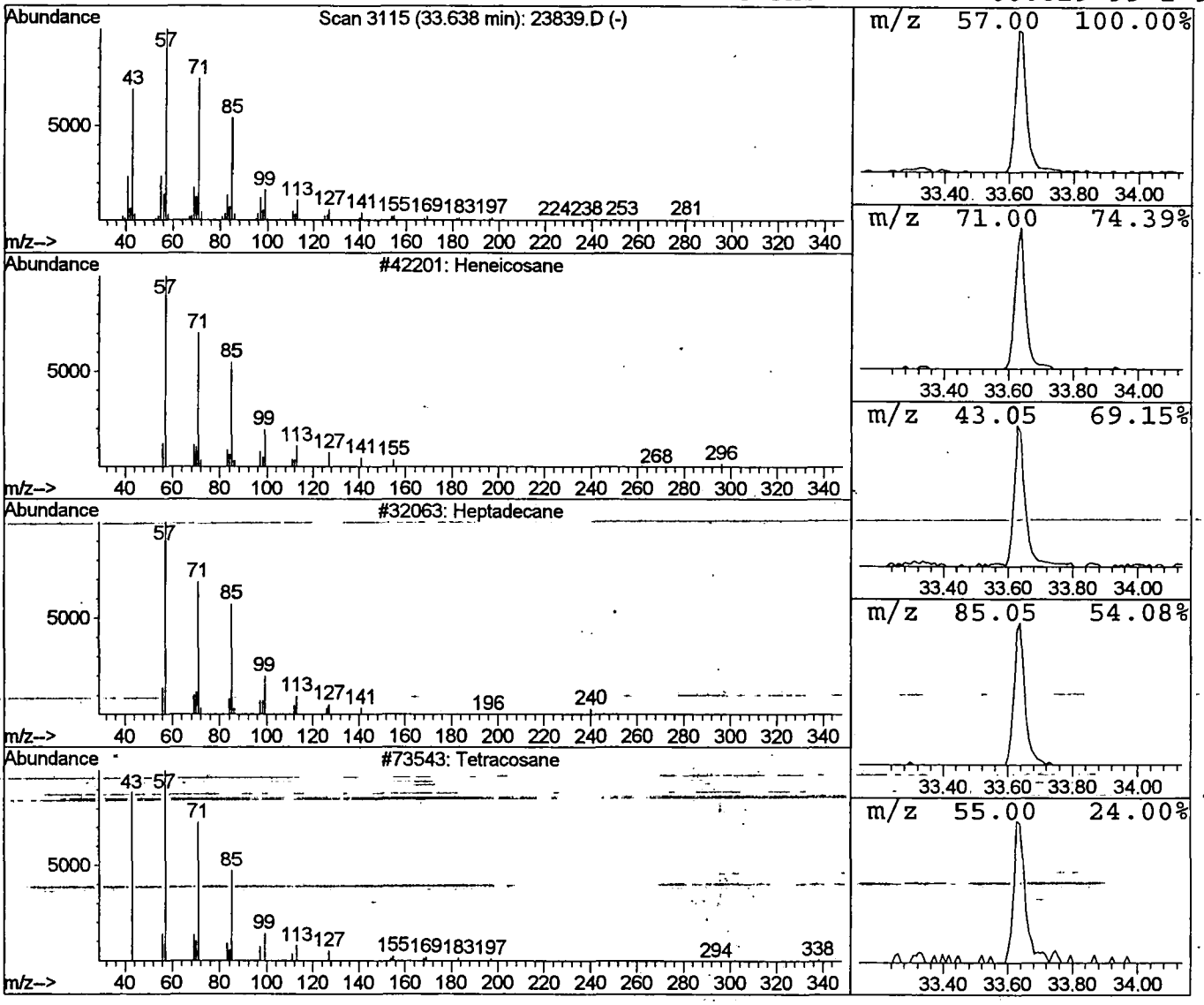
Quant Method : C:\HPCHEM\1\METHODS\NP101901.M (RTE Integrator)
Title : 209 625/TCLP calibration table
Library : C:\DATABASE\NBS75K.L

****

Peak Number 8 Heneicosane Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
33.64	0.77 ug/l	185301	Chrysene-d12	33.06

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heneicosane	296	C21H44	000629-94-7	91
2	Heptadecane	240	C17H36	000629-78-7	91
3	Tetracosane	338	C24H50	000646-31-1	91
4	Pentacosane	352	C25H52	000629-99-2	91



NYC-WTC_000089380

Library Search Compound Report

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Acq On : 31 Oct 2001 2:14 am
Sample : gc/ms unknown
Misc :
MS Integration Params: LSCINT.P

Vial: 13
Operator: 209 GALOS
Inst : GC/MS 209
Multiplr: 1.00

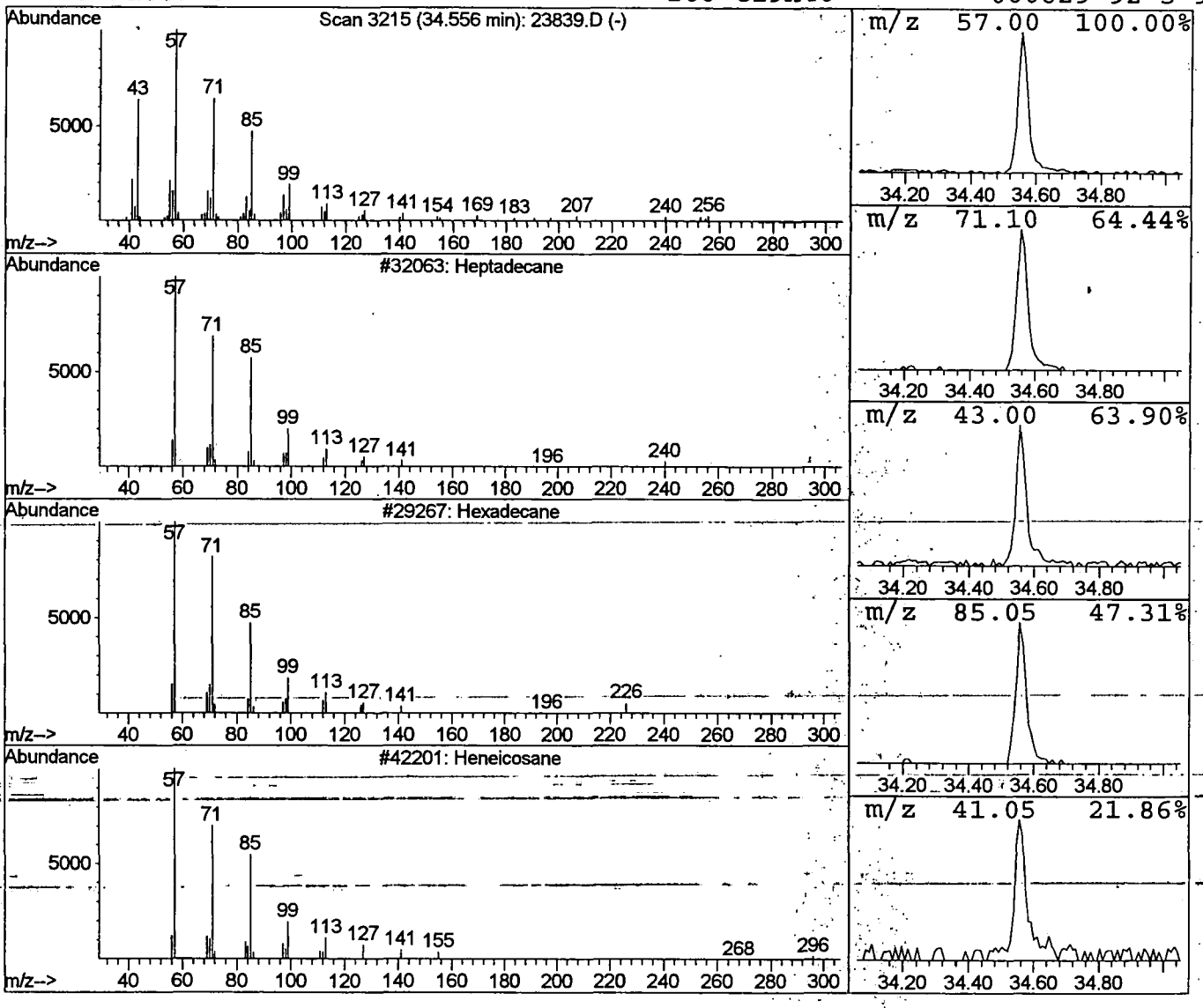
Quant Method : C:\HPCHEM\1\METHODS\NP101901.M (RTE Integrator)
Title : 209 625/TCLP calibration table
Library : C:\DATABASE\NBS75K.L

****

Peak Number 9 Heptadecane — Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
34.56	0.54 ug/l	129695	Chrysene-d12	33.06

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heptadecane	240	C17H36	000629-78-7	95
2	Hexadecane	226	C16H34	000544-76-3	91
3	Heneicosane	296	C21H44	000629-94-7	91
4	Nonadecane	268	C19H40	000629-92-5	91



Library Search Compound Report

Data File : C:\HPCHEM\1\DATA\OCT3001\23839.D
Acq On : 31 Oct 2001 2:14 am
Sample : gc/ms unknown
Misc :
MS Integration Params: LSCINT.P

Vial: 13
Operator: 209 GALOS
Inst : GC/MS 209
Multiplr: 1.00

Quant Method : C:\HPCHEM\1\METHODS\NP101901.M (RTE Integrator)
Title : 209 625/TCLP calibration table
Library : C:\DATABASE\NBS75K.L

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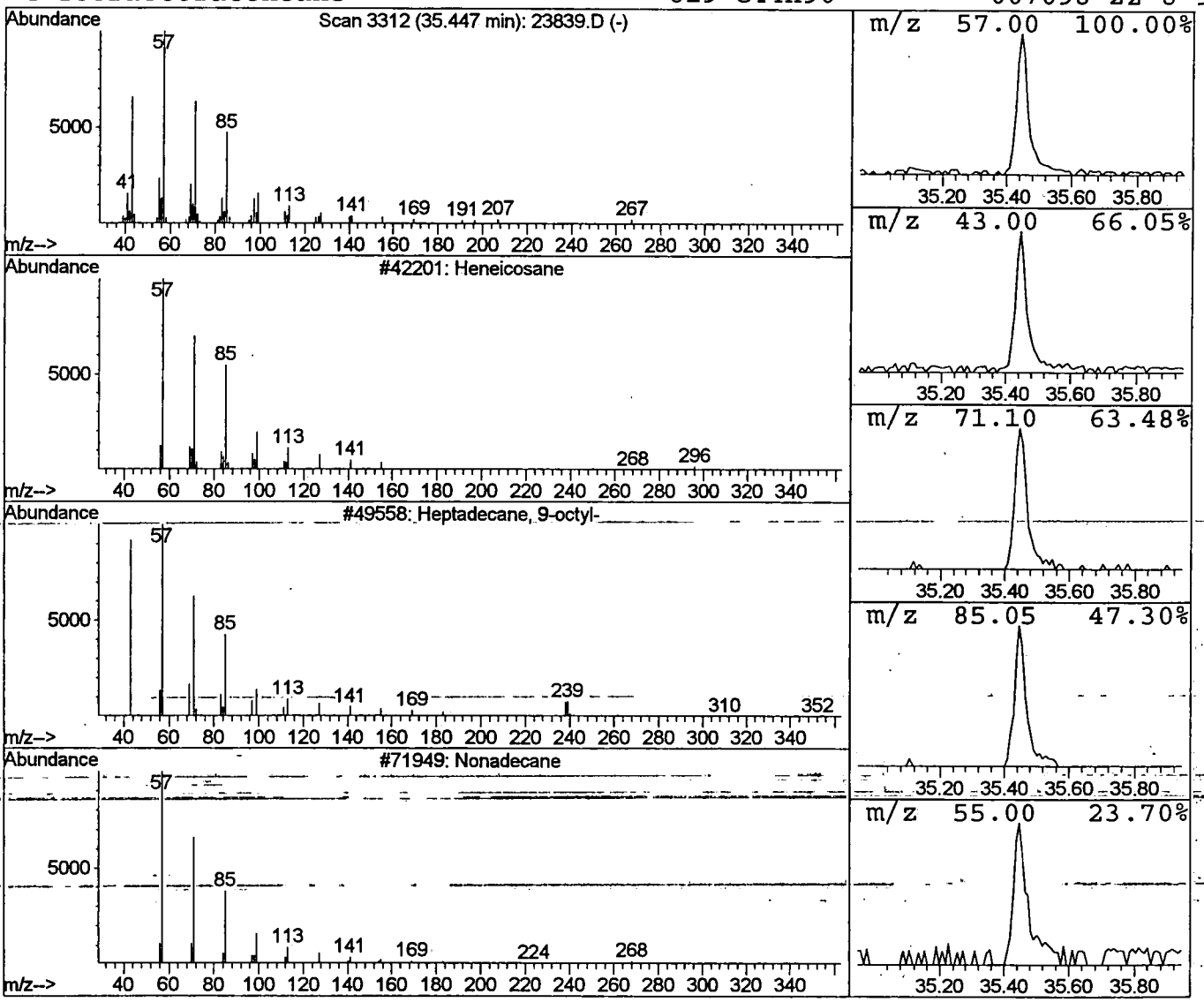
Peak Number 10 Heneicosane Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
35.45	0.56 ug/l	110329	Perylene-d12	37.16

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heneicosane	296	C21H44	000629-94-7	91
2	Heptadecane, 9-octyl-	352	C25H52	007225-64-1	91
3	Nonadecane	268	C19H40	000629-92-5	91
4	Tetratetracontane	619	C44H90	007098-22-8	91



NYC 9/11 Public Portal Document

Tentatively Identified Compound (LSC) summary

Operator ID: 209 GALOS Date Acquired: 31 Oct 2001 2:14 am
Data File: C:\HPCHEM\1\DATA\OCT3001\23839.D
Name: gc/ms unknown
Misc:
Method: C:\HPCHEM\1\METHODS\NP101901.M (RTE Integrator)
Title: 209 625/TCLP calibration table
Library Searched: C:\DATABASE\NBS75K.L

TIC Top Hit name	RT	EstConc	Units	Area	IntStd	ISRT	ISArea	ISConc
2-Hexanone	6.43	0.5	ug/l	124456	ISTD01	11.72	5165660	20.0
2-Pentanone, 4-hydro	7.69	0.6	ug/l	144202	ISTD01	11.72	5165660	20.0
1,4 Dichlorobenzene-	11.57	0.6	ug/l	144313	ISTD01	11.72	5165660	20.0
Butyl hexadecanoate	29.45	1.2	ug/l	291219	ISTD05	33.06	4798890	20.0
Butyl tetradecanoate	31.59	0.9	ug/l	224590	ISTD05	33.06	4798890	20.0
Heptadecane	31.68	0.7	ug/l	178368	ISTD05	33.06	4798890	20.0
Octadecane	32.67	0.7	ug/l	164261	ISTD05	33.06	4798890	20.0
Heneicosane	33.64	0.8	ug/l	185301	ISTD05	33.06	4798890	20.0
Heptadecane	34.56	0.5	ug/l	129695	ISTD05	33.06	4798890	20.0
Heneicosane	35.45	0.6	ug/l	110329	ISTD06	37.16	3933740	20.0

23839.D NP103001.M Thu Nov 01 14:36:11 2001

Comments for Sample AD23839 - Analysis $GCMS

Westchester Dept. of Labs & Research
User: Vinci, David L
Date: 11-05-2001 Time: 15:59:37

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 present in the Laboratory Blank.