User: Galos, Marie
Westchester Dept. of Labs & Research
Date: 11/19/2001 Time: 14:57:34

Sample: AD25038
Analysis: $GCMS GCMS Scan For Unknowns
Units: ug
Started: 11/19/01 14:57 Ended: 11/19/01 14:57 Analyst: MG
Page: 1

	Component Name	Viol	Result
1	Other unknown compounds		see comment

Comments for Sample AD25038 - Analysis $GCMS

Westchester Dept. of Labs & Research
User: Vinci, David L
Date: 11-21-2001 Time: 11:47:34

A scan of the extract prepared for the 508 analysis, from 35-550 amu using a mass spectrometer, does not reveal the presence of non-target compounds, except for siloxanes which are due to column bleed.

Quantitation Report (QT Reviewed)

Data File : C:\HPCHEM\1\DATA\NOV0801\25038.D Vial: 12
Acq On : 9 Nov 2001 1:21 am Operator: 209 GALOS
Sample : gc/ms unknown Inst : GC/MS 209
Misc : Multiplr: 1.00
MS Integration Params: RTEINT.P
Quant Time: Nov 9 2:09 2001 Quant Results File: NP103001.RES

Quant Method : C:\HPCHEM\1\METHODS\NP103001.M (RTE Integrator)
Title : 209 625/TCLP calibration table
Last Update : Fri Nov 02 16:41:16 2001
Response via : Initial Calibration
DataAcq Meth : NP103001

Internal Standards	R.T.	QIon	Response	Conc Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	11.71	152	789872	20.00 ug/l	-0.01
19) Naphthalene-d8	15.32	136	3008698	20.00 ug/l	-0.01
31) Acenaphthene-d10	20.60	164	1485489	20.00 ug/l	-0.01
49) Phenanthrene-d10	25.02	188	2187783	20.00 ug/l	-0.01
59) Chrysene-d12	33.06	240	1593858	20.00 ug/l	-0.01
68) Perylene-d12	37.16	264	1237295	20.00 ug/l	0.00
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	11.70	132	446	0.01 ug/l	0.47
Spiked Amount 75.000			Recovery =	0.01%	
7) 1,2-Dichlorobenzene-d4	12.24	152	9344	0.26 ug/l	-0.01
Spiked Amount 50.000			Recovery =	0.52%	
20) Nitrobenzene-d5	0.00	82	0	0.00 ug/l	
Spiked Amount 50.000			Recovery =	0.00%	
32) 2-Fluorobiphenyl	18.74	172	3245	0.03 ug/l	0.13
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.	
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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	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	15.38	128	668	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	20.93	165	739	N.D.	
45)	Diethylphthalate	22.10	149	335	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	25.02	178	630	N.D.	
56)	Anthracene	25.02	178	630	N.D.	
57)	Di-n-butylphthalate	27.02	149	2426	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	33.06	228	3752	N.D.	
66)	Bis(2-ethylhexyl)phthalate	33.32	149	1016	N.D.	
67)	Chrysene	33.06	228	3752	N.D.	
69)	Di-n-Octylphthalate	0.00	149	0	N.D.	
70)	Benzo(b)fluoranthene	0.00	252	0	N.D.	

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Misc : Multiplr: 1.00
MS Integration Params: RTEINT.P
Quant Time: Nov 9 2:09 2001 Quant Results File: NP103001.RES

Quant Method : C:\HPCHEM\1\METHODS\NP103001.M (RTE Integrator)
Title : 209 625/TCLP calibration table
Last Update : Fri Nov 02 16:41:16 2001
Response via : Initial Calibration
DataAcq Meth : NP103001

	Compound	R.T.	QIon	Response	Conc Unit	Qvalue
71)	Benzo(k)fluoranthene	0.00	252	0	N.D.	
72)	Benzo(a)pyrene	37.16	252	4568	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

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Acq On : 9 Nov 2001 1:21 am
Sample : gc/ms unknown
Misc :
MS Integration Params: RTEINT.P
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Vial: 12
Operator: 209 GALOS
Inst : GC/MS 209
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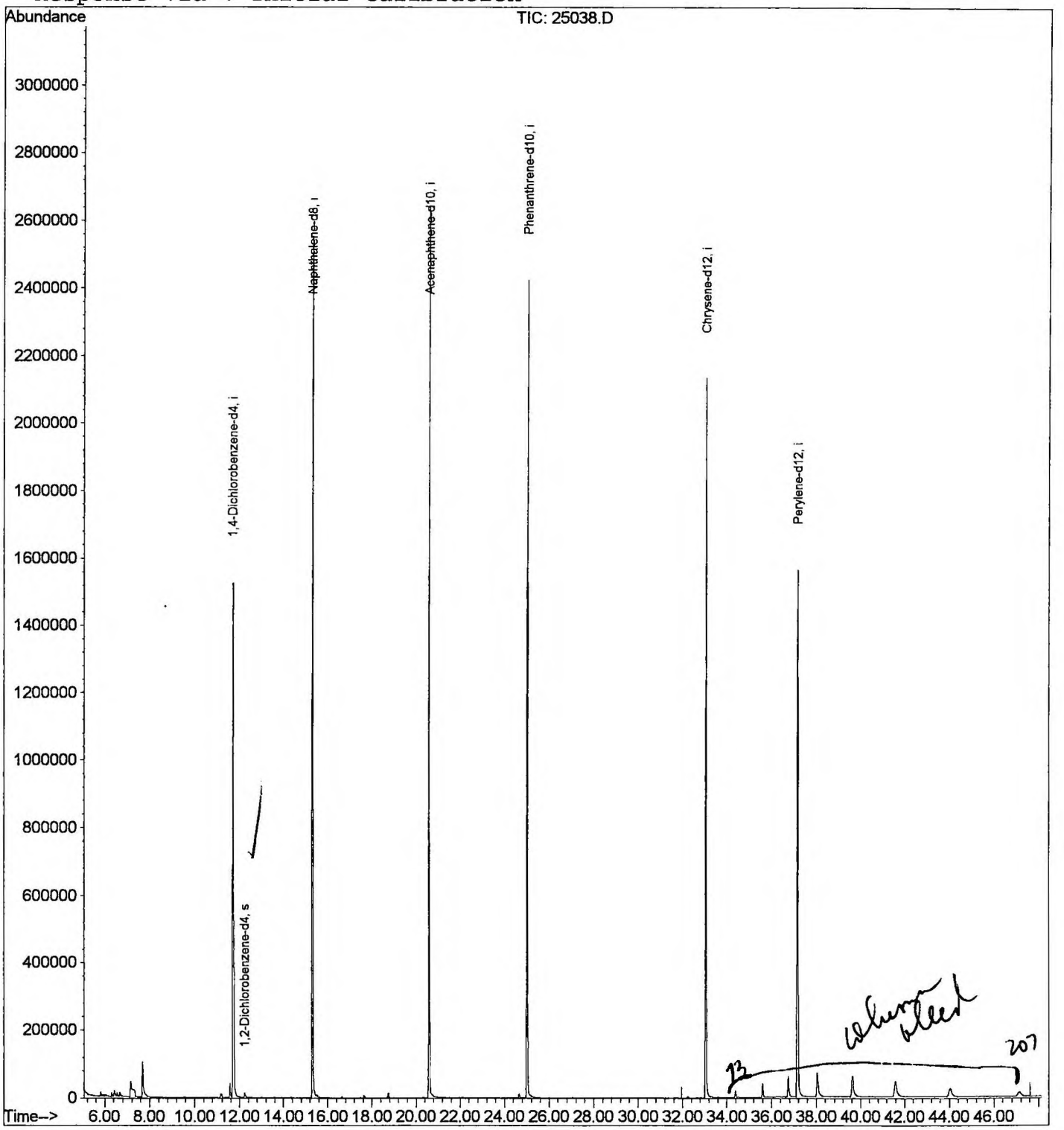
Title : 209 625/TCLP calibration table
Last Update : Fri Nov 02 16:41:16 2001
Response via : Initial Calibration



LSC Area Percent Report

Data File : C:\HPCHEM\1\DATA\NOV0801\25038.D Vial: 12
Acq On : 9 Nov 2001 1:21 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\NP103001.M (RTE Integrator)
Title : 209 625/TCLP calibration table
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

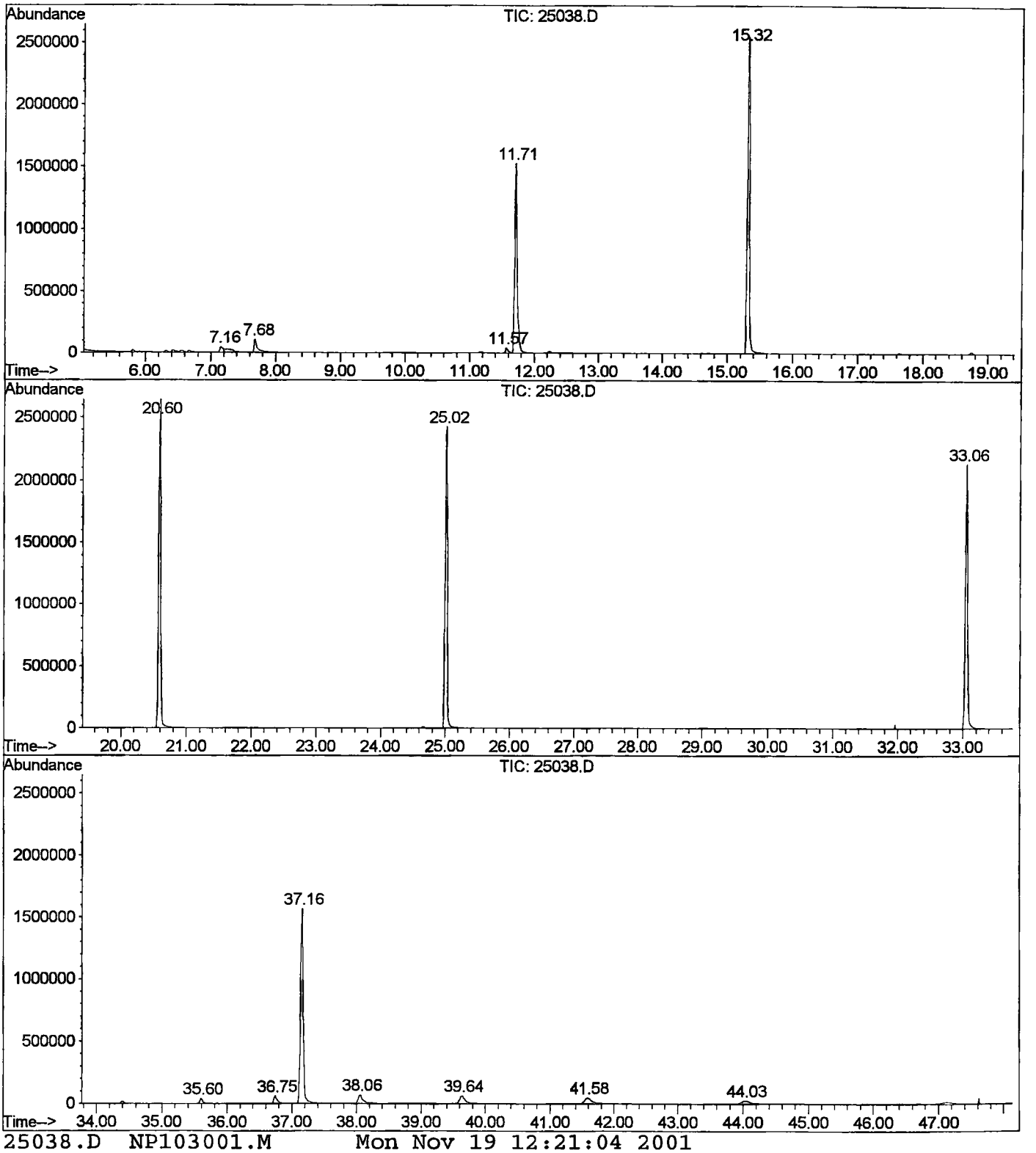
peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	peak area	peak % max.	% of total
1	7.159	224	230	235	rBV	45491	145937	2.49%	0.450%
2	7.682	283	287	312	rBV	102885	328906	5.61%	1.014%
3	11.574	706	711	720	rBV	41675	106531	1.82%	0.329%
4	11.712	720	726	742	rBV	1525813	4063311	69.36%	12.532%
5	15.320	1113	1119	1137	rBV	2551666	5485125	93.64%	16.918%
6	20.599	1687	1694	1717	rBV	2643430	5857951	100.00%	18.067%
7	25.024	2169	2176	2188	rBV2	2423968	5487071	93.67%	16.924%
8	33.056	3044	3051	3068	rBV2	2133525	5005060	85.44%	15.437%
9	35.599	3321	3328	3337	rBV2	40880	114306	1.95%	0.353%
10	36.747	3447	3453	3464	rBV	59372	197429	3.37%	0.609%
11	37.160	3489	3498	3522	rBV	1564054	4484381	76.55%	13.831%
12	38.060	3588	3596	3611	rBV2	68249	305441	5.21%	0.942%
13	39.639	3757	3768	3782	rBV2	60568	337281	5.76%	1.040%
14	41.585	3967	3980	3998	rBV2	45119	314080	5.36%	0.969%
15	44.027	4233	4246	4262	rBV5	23627	189844	3.24%	0.586%

Sum of corrected areas: 32422654

25038.D NP103001.M Mon Nov 19 12:21:01 2001

LSC Report - Integrated Chromatogram

File : C:\HPCHEM\1\DATA\NOV0801\25038.D
Operator : 209 GALOS
Acquired : 9 Nov 2001 1:21 am using AcqMethod NP103001
Instrument : GC/MS 209
Sample Name: gc/ms unknown
Misc Info :
Vial Number: 12
Quant File :NP103001.RES (RTE Integrator)



25038.D NP103001.M Mon Nov 19 12:21:04 2001

Library Search Compound Report

Data File : C:\HPCHEM\1\DATA\NOV0801\25038.D
Acq On : 9 Nov 2001 1:21 am
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\NP103001.M (RTE Integrator)
Title : 209 625/TCLP calibration table
Library : C:\DATABASE\NBS75K.L

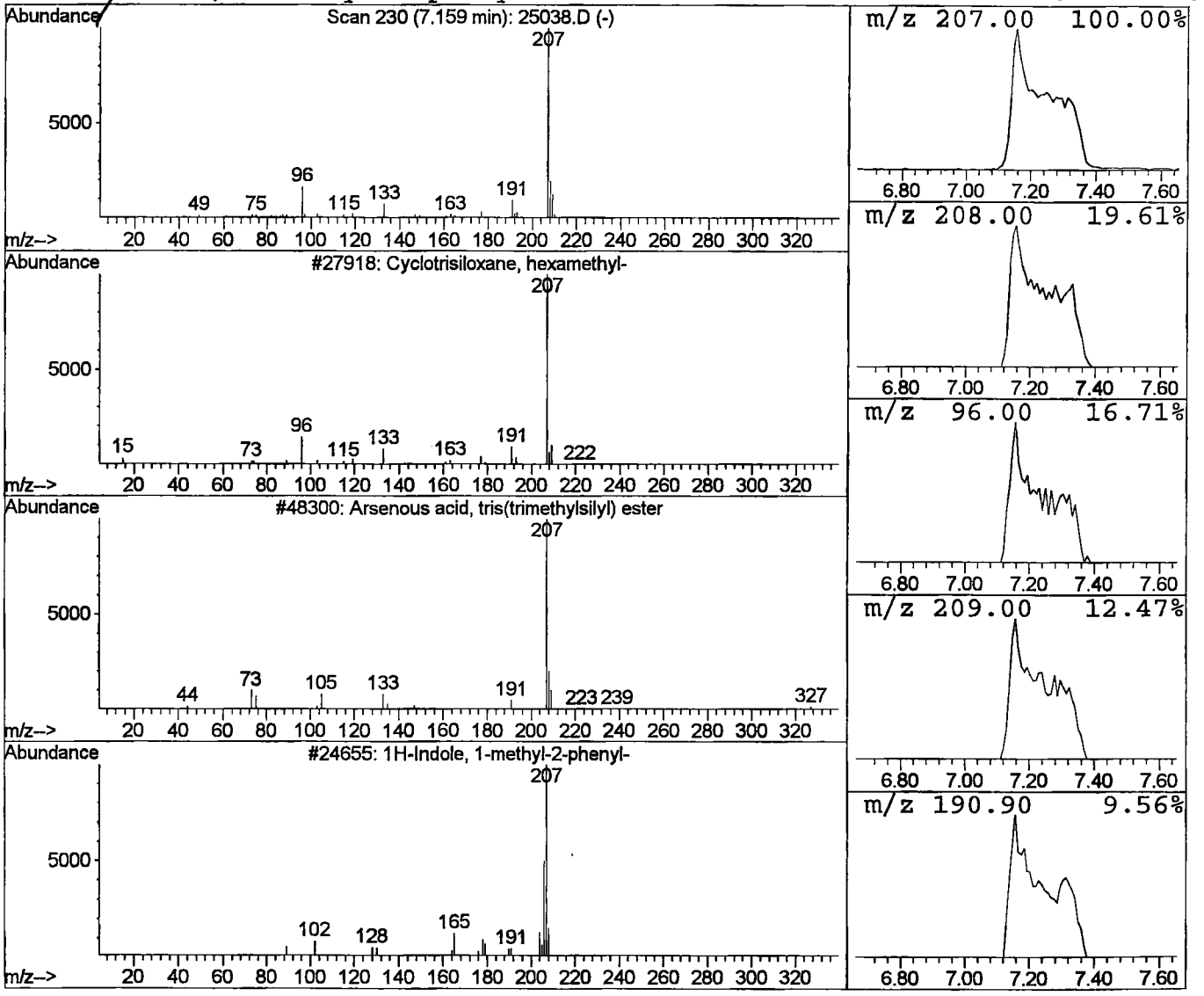
****

Peak Number 1 Cyclotrisiloxane, hexamethyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.16	0.72 ug/l	145937	1,4-Dichlorobenzene-d4	11.71

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclotrisiloxane, hexamethyl-	222	C6H18O3Si3	000541-05-9	80
2	Arsenous acid, tris(trimethylsilyl)	342	C9H27AsO3Si3	055429-29-3	56
3	1H-Indole, 1-methyl-2-phenyl-	207	C15H13N	003558-24-5	39
4	1H-Indole, 2-methyl-3-phenyl-	207	C15H13N	004757-69-1	9

Column



Library Search Compound Report

Data File : C:\HPCHEM\1\DATA\NOV0801\25038.D
Acq On : 9 Nov 2001 1:21 am
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\NP103001.M (RTE Integrator)
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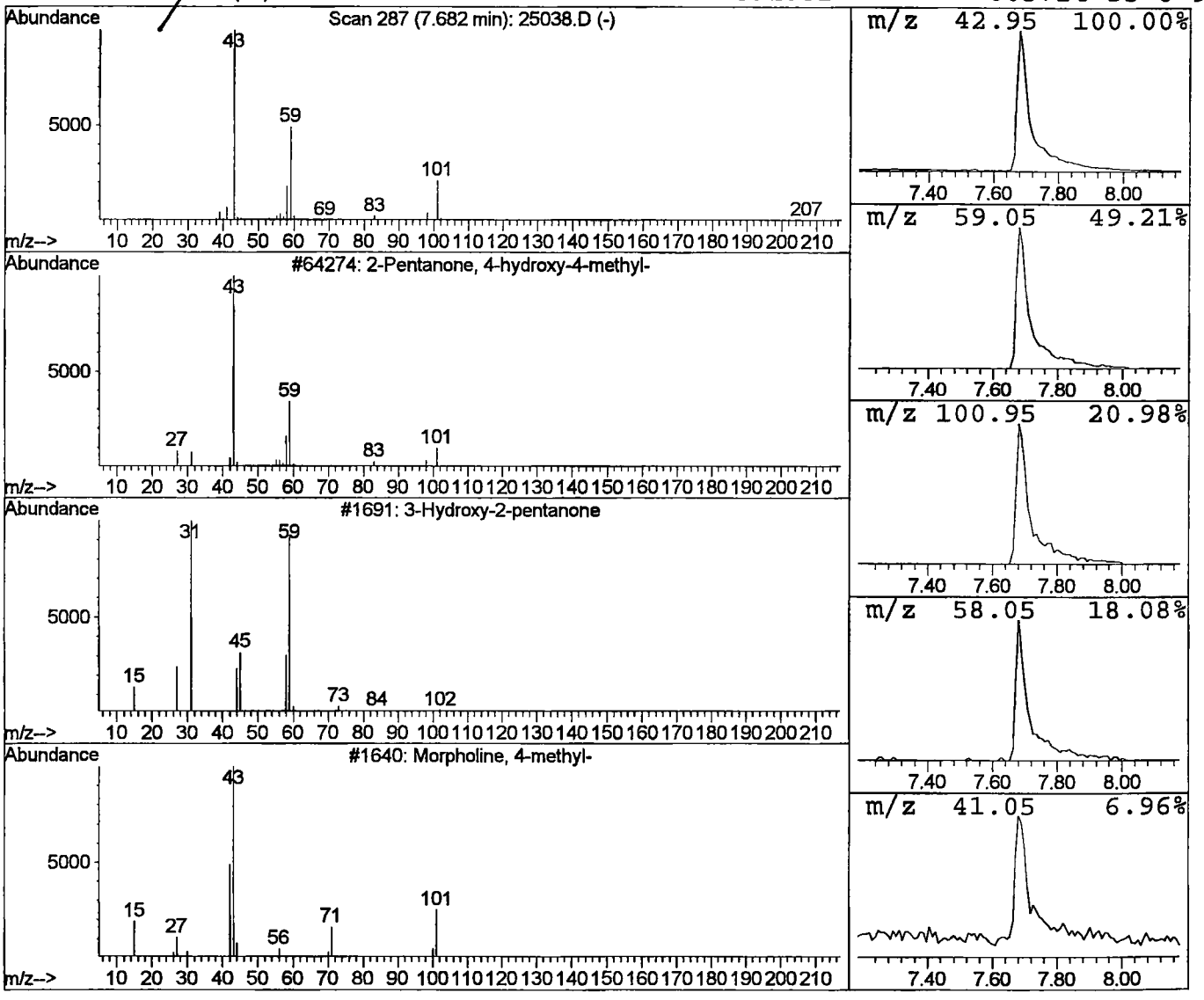
Library : C:\DATABASE\NBS75K.L

****

Peak Number 2 2-Pentanone, 4-hydroxy-4-methy Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.68	1.62 ug/l	328906	1,4-Dichlorobenzene-d4	11.71

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	28
2	3-Hydroxy-2-pentanone	102	C5H10O2	003142-66-3	9
3	Morpholine, 4-methyl-	101	C5H11NO	000109-02-4	9
4	CH2=CHCH2C(O)OCH3	100	C5H8O2	003724-55-8	9



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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\NP103001.M (RTE Integrator)
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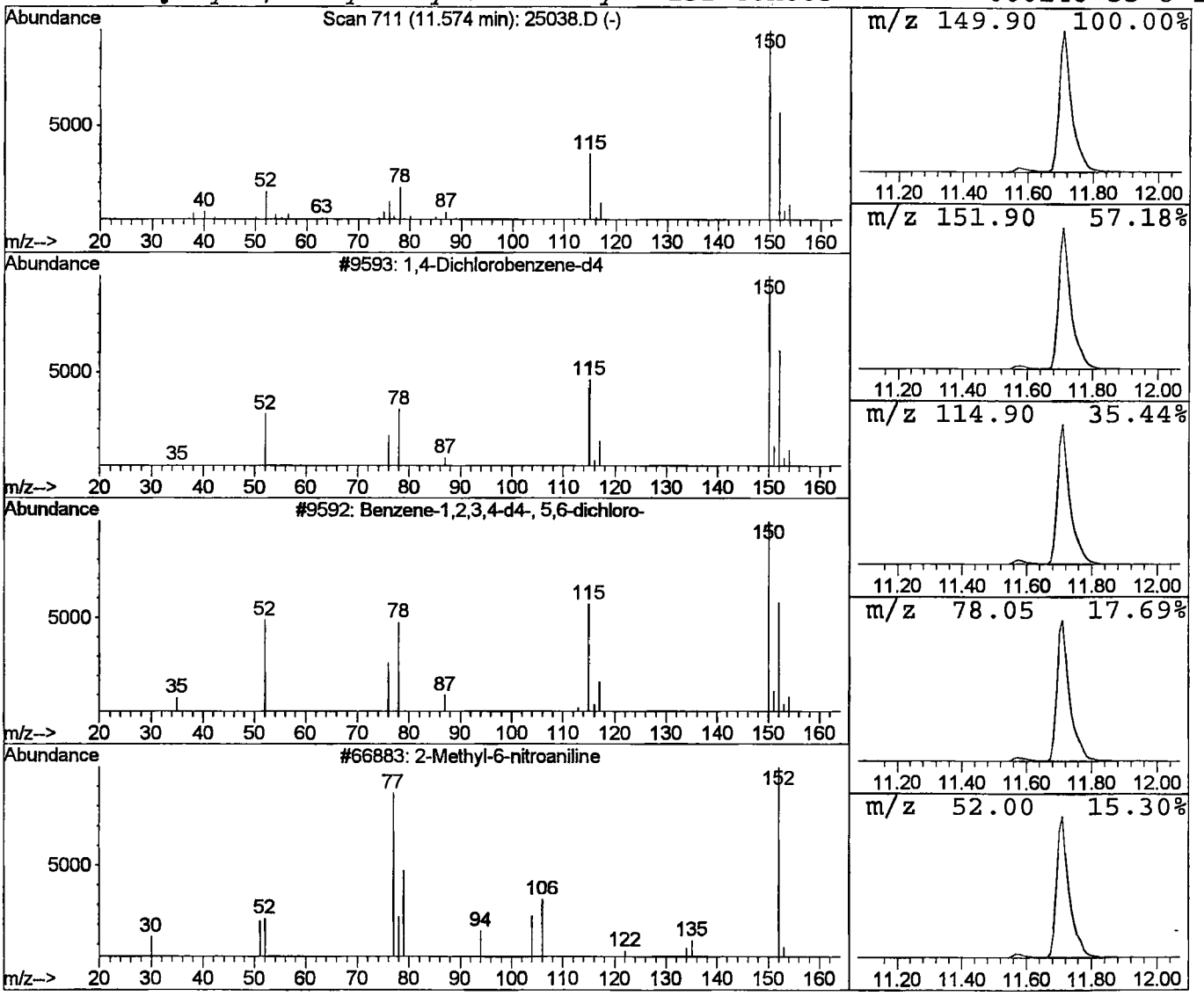
****

Peak Number 3 1,4-Dichlorobenzene-d4 Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.57	0.52 ug/l	106531	1,4-Dichlorobenzene-d4	11.71

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	10
4	Benzaldehyde, 2-hydroxy-3-methoxy-	152	C8H8O3	000148-53-8	10



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Misc :
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Vial: 12
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Multiplr: 1.00

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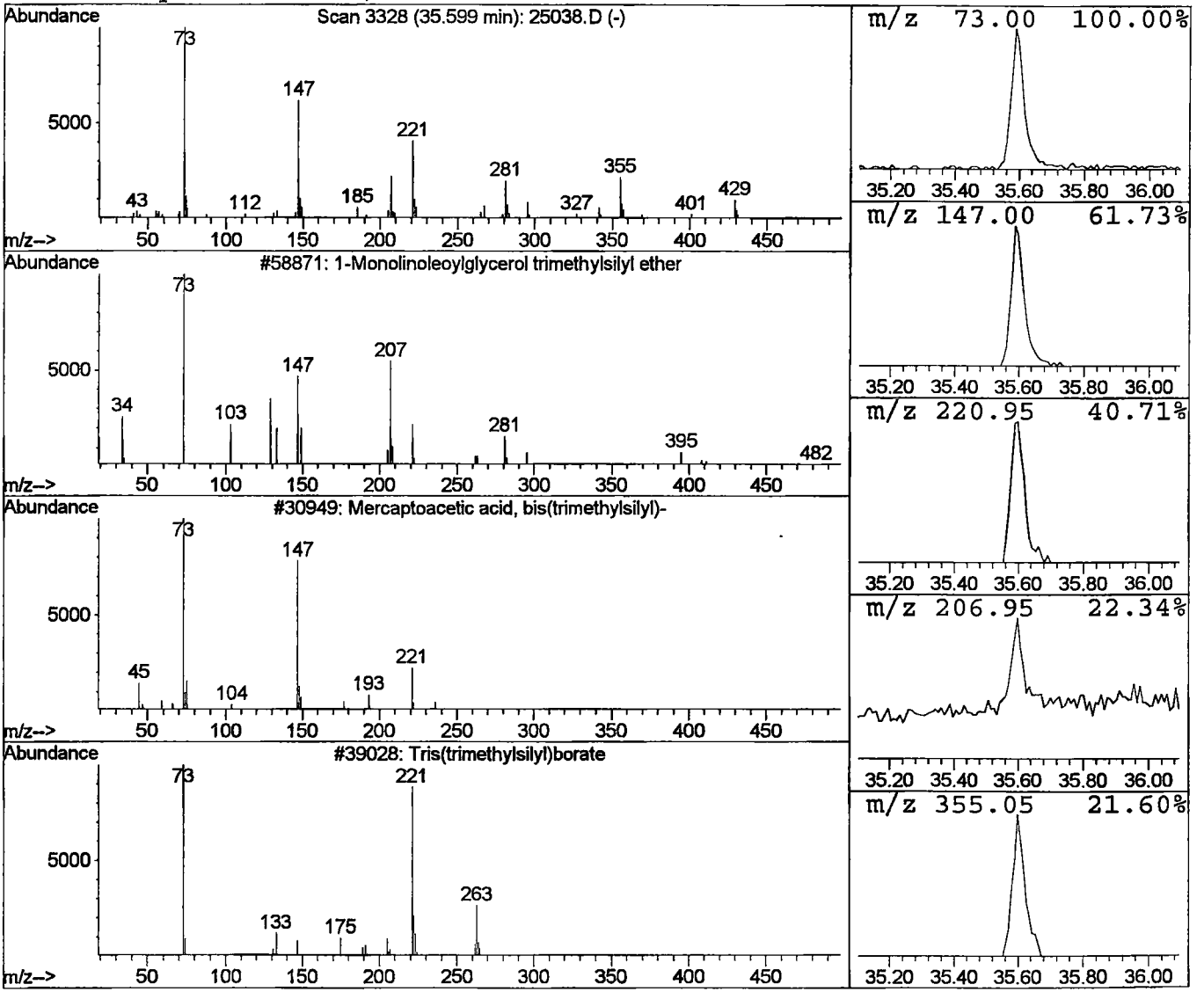
****

Peak Number 4 1-Monolinoleoylglycerol trimet Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
35.60	0.51 ug/l	114306	Perylene-d12	37.16

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Monolinoleoylglycerol trimethylsi	498	C27H54O4Si2	054284-45-6	37
2	Mercaptoacetic acid, bis(trimethyls	236	C8H20O2SSi2	006398-62-5	12
3	Tris(trimethylsilyl)borate	278	C9H27BO3Si3	004325-85-3	10
4	N-Methyladrenaline, tri-TMS	413	C19H39NO3Si3	000000-00-0	9



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Inst : GC/MS 209
Multiplr: 1.00

Quant Method : C:\HPCHEM\1\METHODS\NP103001.M (RTE Integrator)
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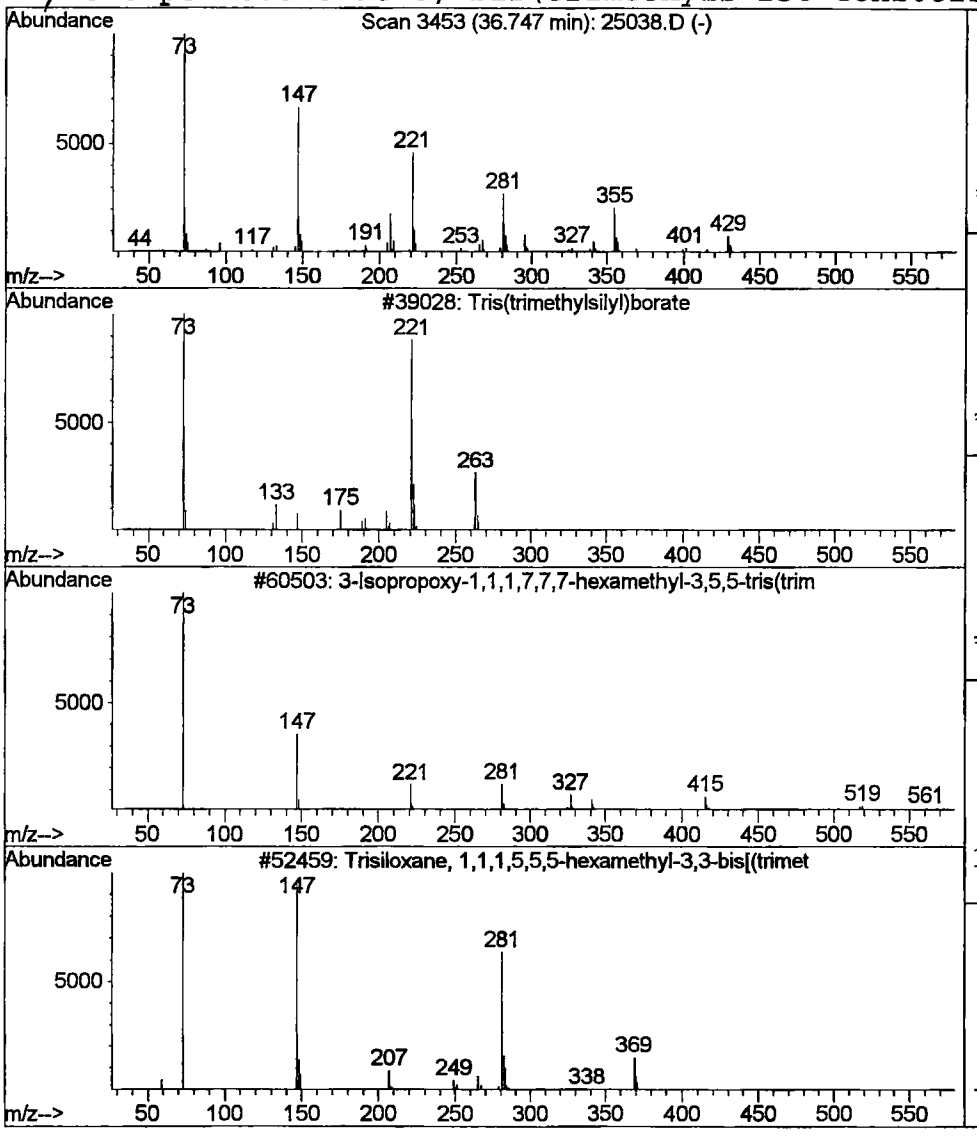
****

Peak Number 5 Tris(trimethylsilyl)borate Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
36.75	0.88 ug/l	197429	Perylene-d12	37.16

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Tris(trimethylsilyl)borate	278	C9H27BO3Si3	004325-85-3	16
2	3-Isopropoxy-1,1,1,7,7,7-hexamethyl	576	C18H52O7Si7	071579-69-6	12
3	Trisiloxane, 1,1,1,5,5,5-hexamethyl	384	C12H36O4Si5	003555-47-3	12
4	Mercaptoacetic acid, bis(trimethyls	236	C8H20O2SSi2	006398-62-5	12



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Misc :
MS Integration Params: LSCINT.P

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

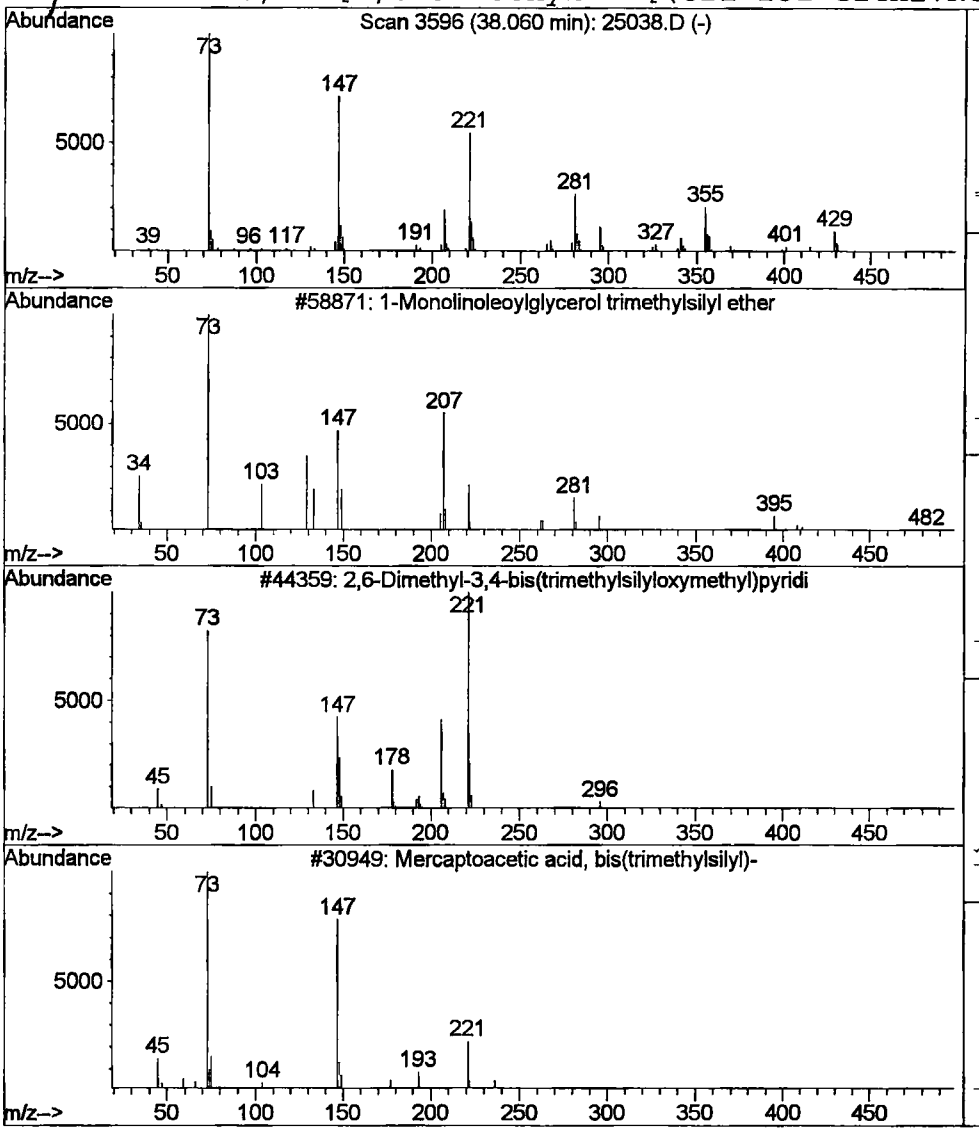
Quant Method : C:\HPCHEM\1\METHODS\NP103001.M (RTE Integrator)
Title : 209 625/TCLP calibration table
Library : C:\DATABASE\NBS75K.L

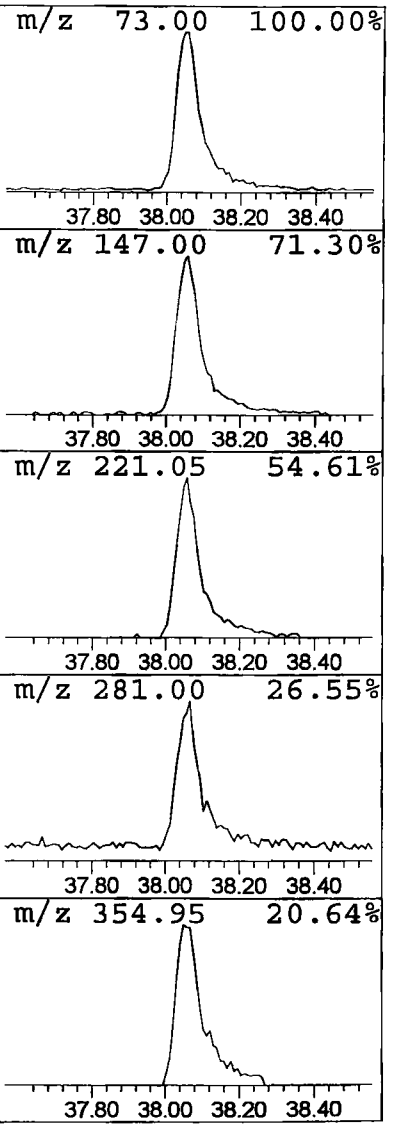
****

Peak Number 6 1-Monolinoleoylglycerol trimet Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
38.06	1.36 ug/l	305441	Perylene-d12	37.16

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Monolinoleoylglycerol trimethylsi	498	C27H54O4Si2	054284-45-6	23
2	2,6-Dimethyl-3,4-bis(trimethylsilyl	311	C15H29NO2Si2	000000-00-0	16
3	Mercaptoacetic acid, bis(trimethyls	236	C8H20O2SSi2	006398-62-5	12
4	Silanamine, N-[2,6-dimethyl-4-[(tri	281	C14H27NOSi2	072088-09-6	11





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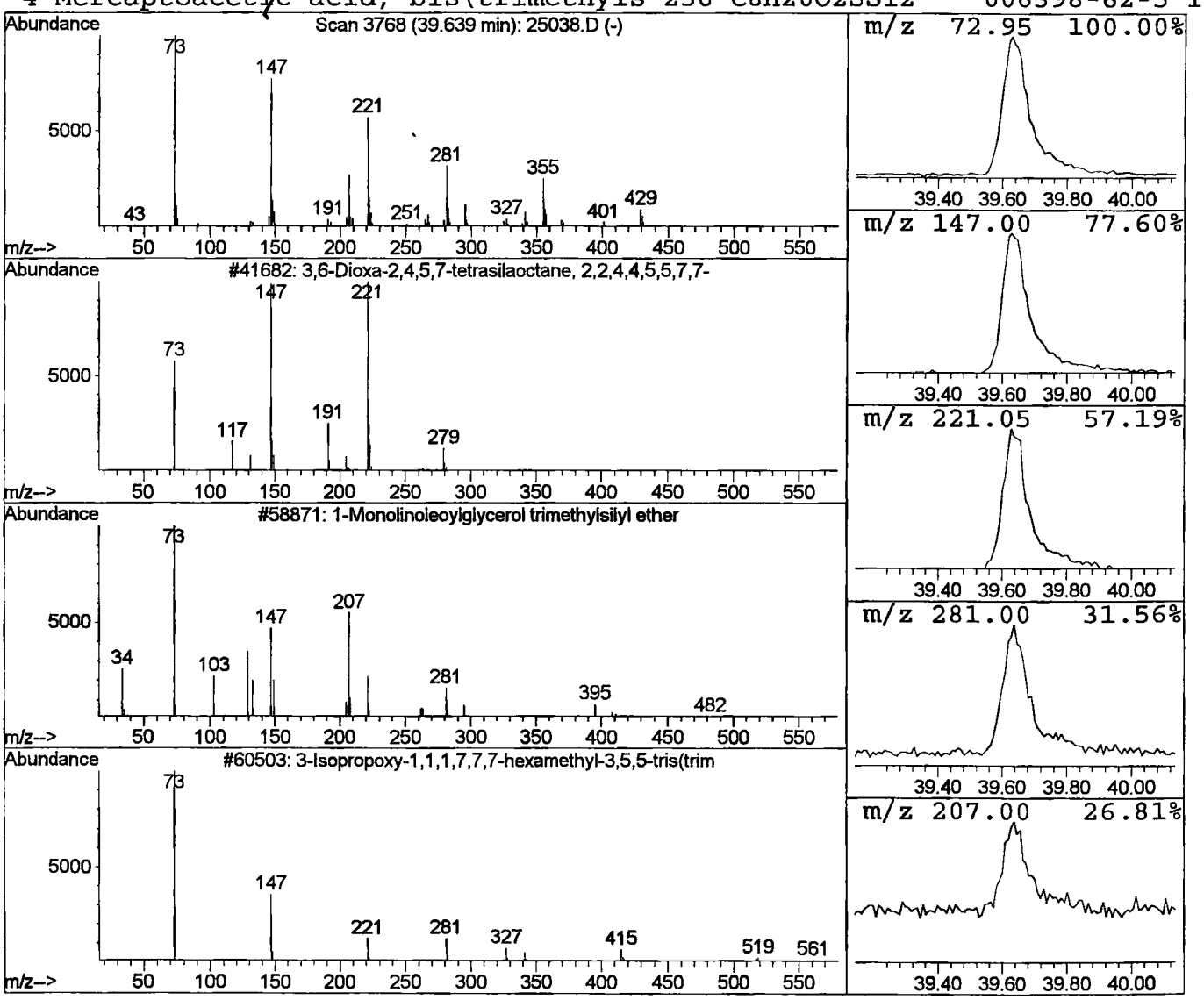
****

Peak Number 7 3,6-Dioxa-2,4,5,7-tetrasilaoct Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
39.64	1.50 ug/l	337281	Perylene-d12	37.16

Column

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	3,6-Dioxa-2,4,5,7-tetrasilaoctane,	294	C10H30O2Si4	004342-25-0	38
2	1-Monolinoleoylglycerol trimethylsi	498	C27H54O4Si2	054284-45-6	33
3	3-Isopropoxy-1,1,1,7,7,7-hexamethyl	576	C18H52O7Si7	071579-69-6	23
4	Mercaptoacetic acid, bis(trimethyls	236	C8H20O2SSi2	006398-62-5	17



Library Search Compound Report

Data File : C:\HPCHEM\1\DATA\NOV0801\25038.D
Acq On : 9 Nov 2001 1:21 am
Sample : gc/ms unknown
Misc :
MS Integration Params: LSCINT.P

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

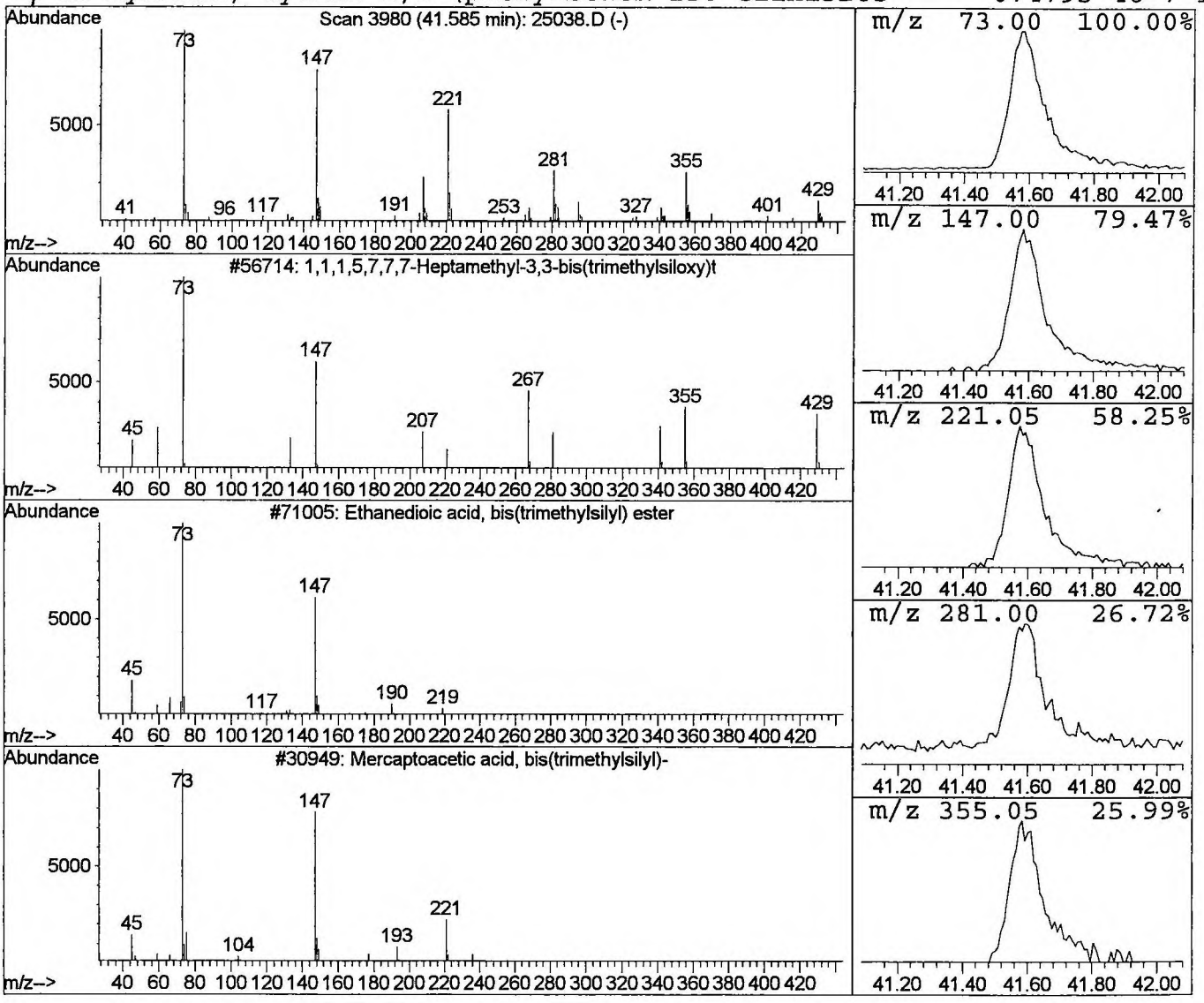
Quant Method : C:\HPCHEM\1\METHODS\NP103001.M (RTE Integrator)
Title : 209 625/TCLP calibration table
Library : C:\DATABASE\NBS75K.L

****

Peak Number 8 1,1,1,5,7,7,7-Heptamethyl-3,3- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
41.58	1.40 ug/l	314080	Perylene-d12	37.16

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,1,1,5,7,7,7-Heptamethyl-3,3-bis(t	444	C13H40O5Si6	038147-00-1	25
2	Ethanedioic acid, bis(trimethylsily	234	C8H18O4Si2	018294-04-7	22
3	Mercaptoacetic acid, bis(trimethyls	236	C8H20O2SSi2	006398-62-5	17
4	dl-Xylitol, cyclic 2,3-(phenylboron	238	C11H15BO5	074793-46-7	12



Library Search Compound Report

Data File : C:\HPCHEM\1\DATA\NOV0801\25038.D
Acq On : 9 Nov 2001 1:21 am
Sample : gc/ms unknown
Misc :
MS Integration Params: LSCINT.P

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

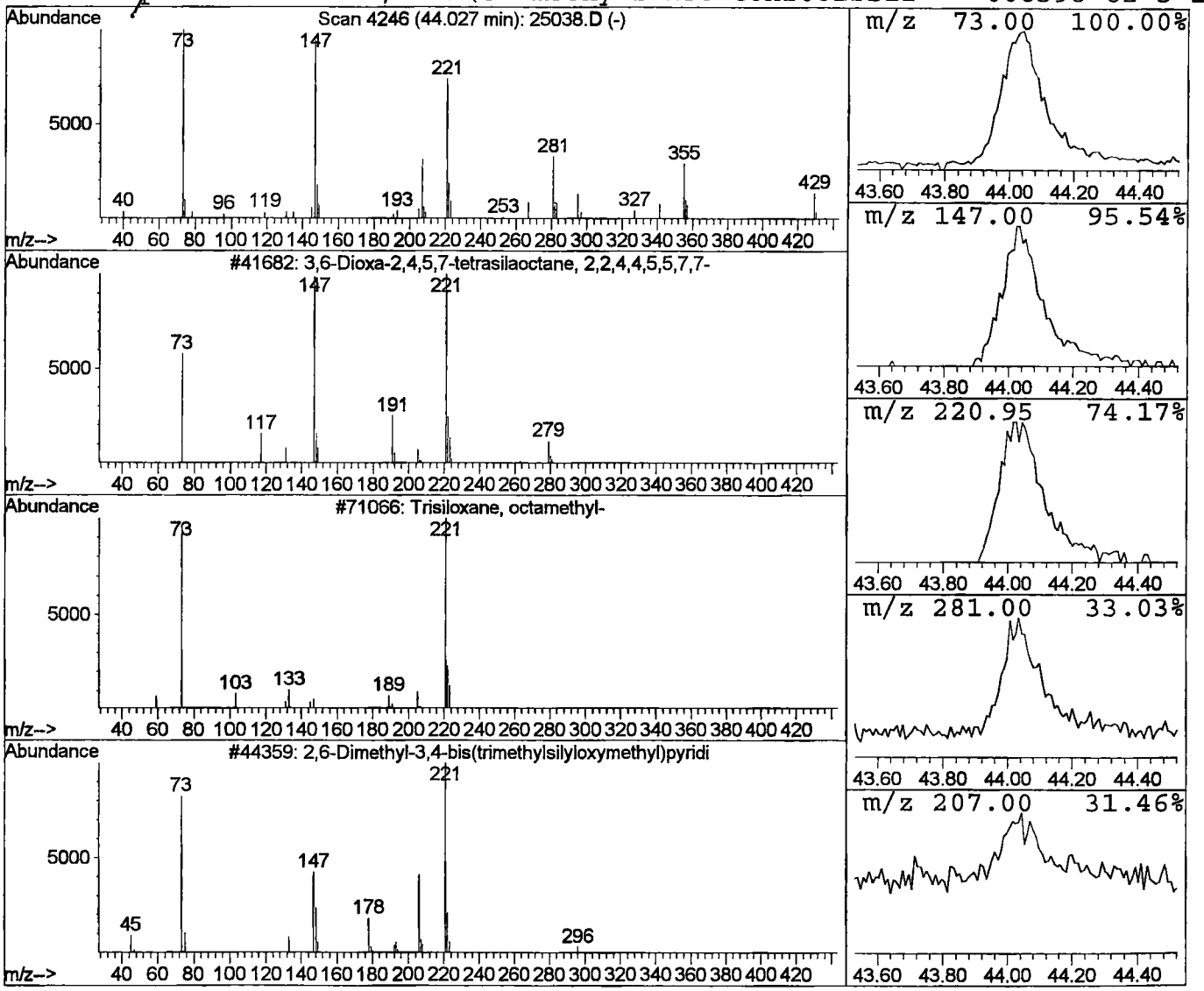
Quant Method : C:\HPCHEM\1\METHODS\NP103001.M (RTE Integrator)
Title : 209 625/TCLP calibration table
Library : C:\DATABASE\NBS75K.L

****

Peak Number 9 3,6-Dioxa-2,4,5,7-tetrasilaoct Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
44.03	0.85 ug/l	189844	Perylene-d12	37.16

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	3,6-Dioxa-2,4,5,7-tetrasilaoctane,	294	C10H30O2Si4	004342-25-0	38
2	Trisiloxane, octamethyl-	236	C8H24O2Si3	000107-51-7	22
3	2,6-Dimethyl-3,4-bis(trimethylsilyl	311	C15H29NO2Si2	000000-00-0	16
4	Mercaptoacetic acid, bis(trimethyls	236	C8H20O2SSi2	006398-62-5	12



Tentatively Identified Compound (LSC) summary

Operator ID: 209 GALOS Date Acquired: 9 Nov 2001 1:21 am
Data File: C:\HPCHEM\1\DATA\NOV0801\25038.D
Name: gc/ms unknown
Misc:
Method: C:\HPCHEM\1\METHODS\NP103001.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
Cyclotrisiloxane, he	7.16	0.7	ug/l	145937	ISTD01	11.71	4063310	20.0
2-Pentanone, 4-hydro	7.68	1.6	ug/l	328906	ISTD01	11.71	4063310	20.0
1,4-Dichlorobenzene-	11.57	0.5	ug/l	106531	ISTD01	11.71	4063310	20.0
1-Monolinoleoylglyce	35.60	0.5	ug/l	114306	ISTD06	37.16	4484380	20.0
Tris(trimethylsilyl)	36.75	0.9	ug/l	197429	ISTD06	37.16	4484380	20.0
1-Monolinoleoylglyce	38.06	1.4	ug/l	305441	ISTD06	37.16	4484380	20.0
3,6-Dioxa-2,4,5,7-te	39.64	1.5	ug/l	337281	ISTD06	37.16	4484380	20.0
1,1,1,5,7,7,7-Heptam	41.58	1.4	ug/l	314080	ISTD06	37.16	4484380	20.0
3,6-Dioxa-2,4,5,7-te	44.03	0.8	ug/l	189844	ISTD06	37.16	4484380	20.0

25038.D NP103001.M Mon Nov 19 12:21:27 2001