User: Vinci, David L
Westchester Dept. of Labs & Research
Date: 01/09/2002 Time: 15:27:40

Sample: AD28143
Analysis: $B1GCMS Blank for GCMS Scan
Units: ug
Started: 11/28/01 15:56 Ended: 11/28/01 15:56 Analyst: DLV
Result source: manual entry
Page: 1

	Component Name	Viol	Result	Qualifier	MDL
1	n-Nitrosodimethylamine		< MDL		2.0
2	bis(2-Chloroethyl)ether		< MDL		2.0
3	1,3-Dichlorobenzene		< MDL		2.0
4	1,4-Dichlorobenzene		< MDL		2.0
5	1,2-Dichlorobenzene		< MDL		2.0
6	2,2'-oxybis(1-Chloropropane)		< MDL		2.0
7	n-Nitrosodi-n-propylamine		< MDL		2.0
8	Hexachloroethane		< MDL		2.0
9	Nitrobenzene		< MDL		2.0
10	Isophorone		< MDL		2.0
11	bis(2-Chloroethoxy)methane		< MDL		2.0
12	1,2,4-Trichlorobenzene		< MDL		2.0
13	Naphthalene		< MDL		2.0
14	Hexachlorobutadiene		< MDL		2.0
15	Hexachlorocyclopentadiene		< MDL		2.0
16	2-Chloronaphthalene		< MDL		2.0
17	Dimethylphthalate		< MDL		2.0
18	2,6-Dinitrotoluene		< MDL		2.0
19	Acenaphthylene		< MDL		2.0
20	Acenaphthene		< MDL		2.0
21	2,4-Dinitrotoluene		< MDL		2.0
22	Diethylphthalate		< MDL		2.0
23	4-Chlorophenylphenylether		< MDL		2.0
24	Fluorene		< MDL		2.0
25	Azobenzene/1,2-Diphenylhydra		< MDL		2.0
26	n-Nitrosodiphenylamine		< MDL		2.0
27	4-Bromophenylphenylether		< MDL		2.0
28	Hexachlorobenzene		< MDL		2.0
29	Phenanthrene		< MDL		2.0
30	Anthracene		< MDL		2.0
31	Di-n-butylphthalate		< MDL		2.0
32	Fluoranthene		< MDL		2.0
33	Pyrene		< MDL		2.0
34	Butylbenzylphthalate		< MDL		2.0
35	Benzo(a)anthracene		< MDL		2.0
36	bis(2-Ethylhexyl)phthalate		< MDL		2.0
37	Chrysene		< MDL		2.0
38	Di-n-octylphthalate		< MDL		2.0
39	Benzo(b)fluoranthene		< MDL		2.0
40	Benzo(k)fluoranthene		< MDL		2.0
41	Benzo(a)pyrene		< MDL		2.0
42	Indeno(1,2,3-cd)pyrene		< MDL		2.0
43	Dibenz(a,h)anthracene		< MDL		2.0
44	Benzo(g,h,i)perylene		< MDL		2.0

Quantitation Report (QT Reviewed)

Data File : M:\INSTRU~1\209\DATA\NOV2601\LBA.D Vial: 9
Acq On : 26 Nov 2001 6:31 pm Operator: 209 GALOS
Sample : 11/20 lab blk scan Inst : GC/MS 209
Misc : Multiplr: 1.00
MS Integration Params: RTEINT.P
Quant Time: Nov 27 12:32 2001 Quant Results File: GCMS1126.RES

Quant Method : C:\HPCHEM\1\METHODS\GCMS1126.M (RTE Integrator)
Title : 209 625/TCLP calibration table
Last Update : Tue Nov 27 12:20:14 2001
Response via : Initial Calibration
DataAcq Meth : NP103001

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	11.70	152	976317	20.00	ug/l	0.00
15) Naphthalene-d8	15.31	136	3744805	20.00	ug/l	0.00
26) Acenaphthene-d10	20.57	164	1840757	20.00	ug/l	0.00
42) Phenanthrene-d10	24.99	188	2644443	20.00	ug/l	0.00
53) Chrysene-d12	33.02	240	1495796	20.00	ug/l	-0.01
59) Perylene-d12	37.12	264	1076366	20.00	ug/l	0.00
System Monitoring Compounds						
52) 2,4-DDD	30.07	235	106226	3.37	ug/mL	0.00
Spiked Amount 5.000			Recovery = 67.40%			
Target Compounds						Qvalue
2) Pyridine	0.00	79	0	N.D.		
3) N-nitroso-dimethyl amine	0.00	74	0	N.D.		
4) Phenol	11.07	94	478	N.D.		
5) bis(2-Chloroethyl)ether	0.00	93	0	N.D.		
6) 2-Chlorophenol	0.00	128	0	N.D.		
7) 1,3-Dichlorobenzene	0.00	146	0	N.D.		
8) 1,4-Dichlorobenzene	0.00	146	0	N.D.		
9) 1,2-Dichlorobenzene	0.00	146	0	N.D.		
10) 2-Methylphenol	0.00	108	0	N.D.		
11) 2,2'-oxybis(1-Chloropropan	0.00	45	0	N.D.		
12) 3&4-Methylphenol	0.00	108	0	N.D.		
13) N-Nitroso-di-n-propylamine	0.00	70	0	N.D.		
14) Hexachloroethane	0.00	117	0	N.D.		
16) Nitrobenzene	0.00	77	0	N.D.		
17) Isophorone	0.00	82	0	N.D.		
18) 2-Nitrophenol	0.00	139	0	N.D.		
19) 2,4-Dimethylphenol	0.00	107	0	N.D.		
20) bis(2-Chloroethoxy)methane	0.00	93	0	N.D.		
21) 2,4-Dichlorophenol	0.00	162	0	N.D.		
22) 1,2,4-Trichlorobenzene	0.00	180	0	N.D.		
23) Naphthalene	15.36	128	865	N.D.		
24) Hexachlorobutadiene	0.00	225	0	N.D.		
25) 4-Chloro-3-methylphenol	0.00	107	0	N.D.		
27) Hexachlorocyclopentadiene	0.00	237	0	N.D.		
28) 2,4,6-Trichlorophenol	0.00	196	0	N.D.		

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

Quantitation Report (QT Reviewed)

Data File : M:\INSTRU~1\209\DATA\NOV2601\LBA.D Vial: 9
Acq On : 26 Nov 2001 6:31 pm Operator: 209 GALOS
Sample : 11/20 lab blk scan Inst : GC/MS 209
Misc : Multiplr: 1.00
MS Integration Params: RTEINT.P
Quant Time: Nov 27 12:32 2001 Quant Results File: GCMS1126.RES

Quant Method : C:\HPCHEM\1\METHODS\GCMS1126.M (RTE Integrator)
Title : 209 625/TCLP calibration table
Last Update : Tue Nov 27 12:20:14 2001
Response via : Initial Calibration
DataAcq Meth : NP103001

	Compound	R.T.	QIon	Response	Conc Unit	Qvalue
29)	2,4,5-Trichlorophenol	0.00	196	0	N.D.	
30)	2-Chloronaphthalene	0.00	162	0	N.D.	
31)	Dimethylphthalate	19.99	163	355	N.D.	
32)	2,6-Dinitrotoluene	0.00	165	0	N.D.	
33)	Acenaphthylene	0.00	152	0	N.D.	
34)	Acenaphthene	0.00	153	0	N.D.	
35)	2,4-Dinitrophenol	0.00	184	0	N.D.	
36)	4-Nitrophenol	0.00	65	0	N.D.	
37)	2,4-Dinitrotoluene	21.03	165	680	N.D.	
38)	Diethylphthalate	22.09	149	1993	N.D.	
39)	4-Chlorophenyl-phenylether	0.00	204	0	N.D.	
40)	Fluorene	0.00	166	0	N.D.	
41)	Azobenzen/ 1,2-Diphenylhyd	22.70	77	199	N.D.	
43)	4,6-Dinitro-2-methylphenol	0.00	198	0	N.D.	
44)	N-Nitrosodiphenylamine	0.00	168	0	N.D.	
45)	4-Bromophenyl-phenylether	0.00	248	0	N.D.	
46)	Hexachlorobenzene	0.00	284	0	N.D.	
47)	Pentachlorophenol	0.00	266	0	N.D.	
48)	Phenanthrene	25.05	178	1478	N.D.	
49)	Anthracene	25.20	178	1044	N.D.	
50)	Di-n-butylphthalate	27.00	149	15058	N.D.	
51)	Fluoranthene	28.68	202	2612	N.D.	
54)	Pyrene	29.35	202	2704	N.D.	
55)	Butylbenzylphthalate	31.51	149	9516	N.D.	
56)	Benzo(a)anthracene	32.98	228	15087	N.D.	
57)	Bis(2-ethylhexyl)phthalate	33.31	149	54976	0.60 ug/l	98
58)	Chrysene	33.10	228	15990	N.D.	
60)	Di-n-Octylphthalate	35.04	149	37796	0.26 ug/l #	88
61)	Benzo(b)fluoranthene	36.06	252	12301	N.D.	
62)	Benzo(k)fluoranthene	36.06	252	10873	N.D.	
63)	Benzo(a)pyrene	36.95	252	9986	N.D.	
64)	Indeno(1,2,3-cd)pyrene	40.84	276	5584	N.D.	
65)	Dibenz(a,h)anthracene	40.89	278	3690	N.D.	
66)	Benzo(g,h,i)perylene	41.91	276	7192	N.D.	

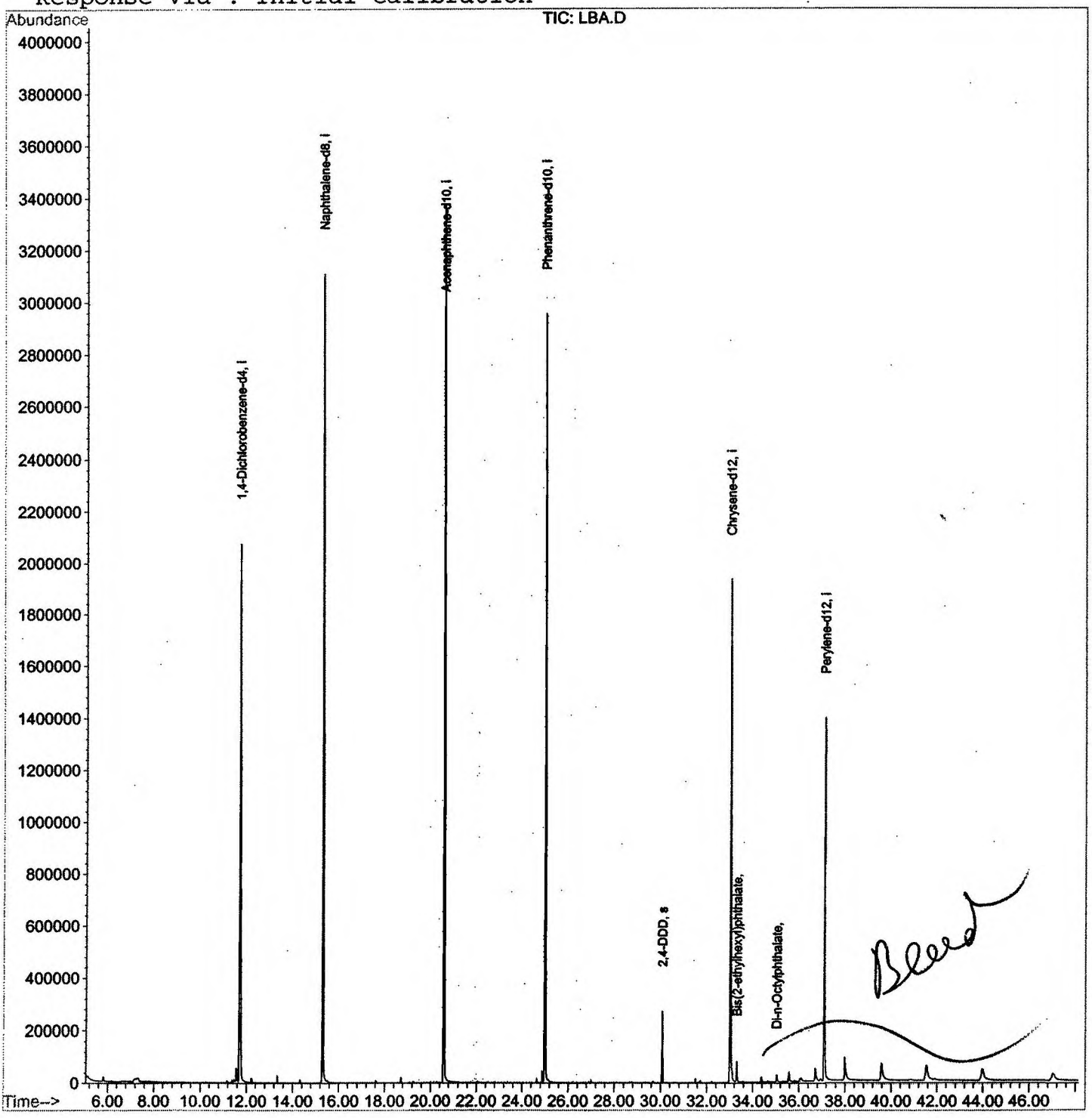
(#) = qualifier out of range (m) = manual integration

Quantitation Report

Data File : M:\INSTRU~1\209\DATA\NOV2601\LBA.D
Acq On : 26 Nov 2001 6:31 pm
Sample : 11/20 lab blk scan
Misc :
MS Integration Params: RTEINT.P
Quant Time: Nov 27 12:32 2001

Vial: 9
Operator: 209 GALOS
Inst : GC/MS 209
Multiplr: 1.00
Quant Results File: GCMS1126.RES

Method : C:\HPCHEM\1\METHODS\GCMS1126.M (RTE Integrator)
Title : 209 625/TCLP calibration table
Last Update : Fri Dec 28 15:11:00 2001
Response via : Initial Calibration



LSC Area Percent Report

Data File : C:\HPCHEM\1\DATA\NOV2601\LBA.D Vial: 9
Acq On : 26 Nov 2001 6:31 pm Operator: 209 GALOS
Sample : 11/20 lab blk scan Inst : GC/MS 209
Misc : Multiplr: 1.00
MS Integration Params: LSCINT.P

Method : C:\HPCHEM\1\METHODS\GCMS1126.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

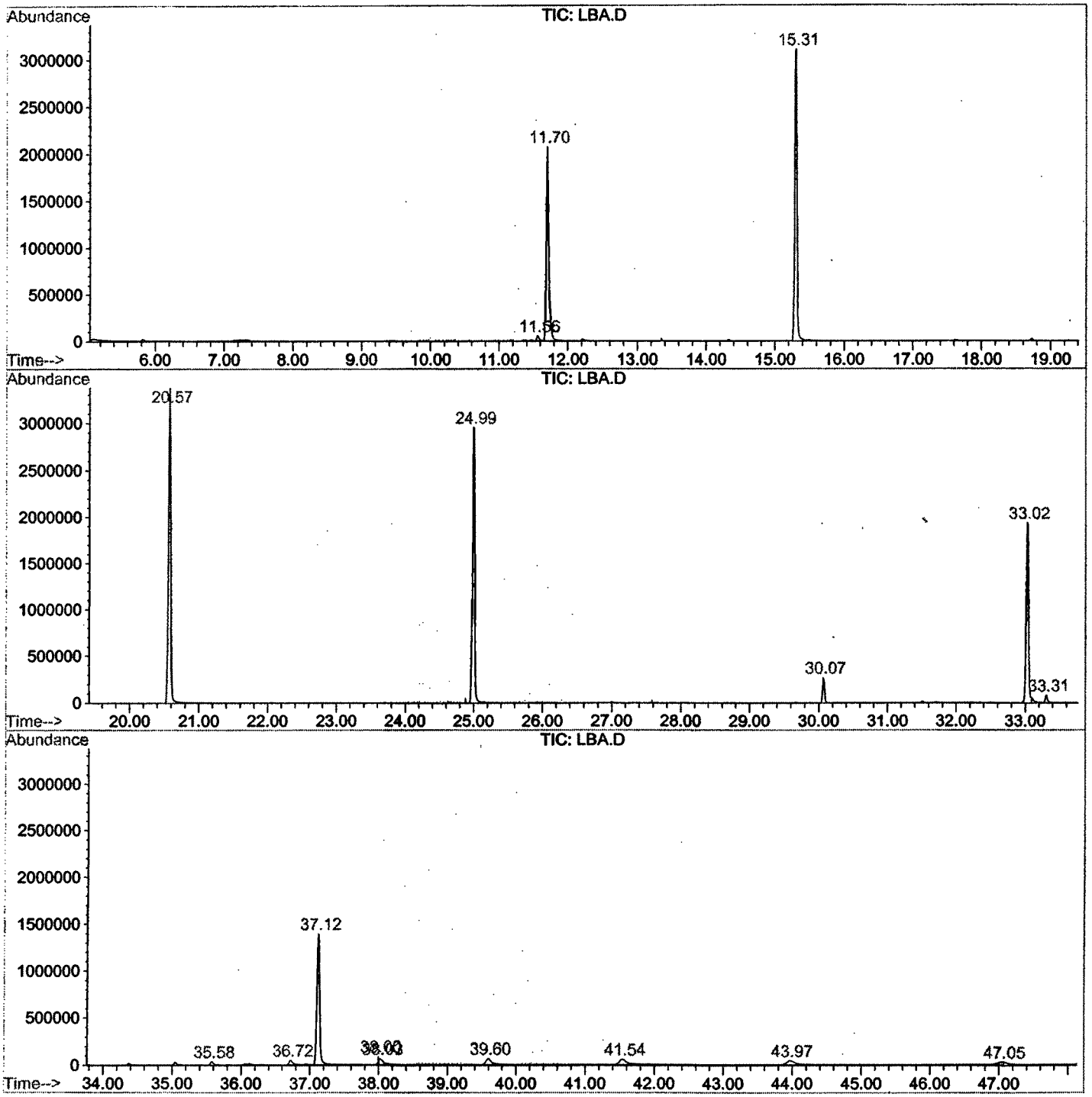
peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	peak area	peak % max.	% of total
1	11.560	706	710	720	rVB	53361	129074	1.79%	0.349%
2	11.697	720	725	743	rBV	2072266	5101607	70.81%	13.780%
3	15.305	1112	1118	1136	rBV	3109673	6811636	94.54%	18.398%
4	20.575	1686	1692	1709	rBV	3376468	7204699	100.00%	19.460%
5	24.991	2167	2173	2186	rBV2	2955565	6615516	91.82%	17.869%
6	30.067	2721	2726	2733	rBV	270994	546992	7.59%	1.477%
7	33.023	3035	3048	3055	rBV2	1934876	4678604	64.94%	12.637%
8	33.308	3072	3079	3086	rBV	79137	176847	2.45%	0.478%
9	35.575	3320	3326	3335	rBV2	37014	115211	1.60%	0.311%
10	36.723	3445	3451	3468	rBV	48439	176851	2.45%	0.478%
11	37.118	3486	3494	3515	rBV2	1390224	3876036	53.80%	10.469%
12	37.999	3585	3590	3591	rBV	87030	85848	1.19%	0.232%
13	38.027	3591	3593	3613	rVB2	60967	219627	3.05%	0.593%
14	39.596	3754	3764	3784	rBV2	66667	401611	5.57%	1.085%
15	41.543	3964	3976	3988	rBV4	57392	383337	5.32%	1.035%
16	43.966	4226	4240	4260	rBV4	42864	365657	5.08%	0.988%
17	47.051	4562	4576	4578	rBV6	24570	133756	1.86%	0.361%

Sum of corrected areas: 37022909

LBA.D GCMS1126.M Thu Jan 03 12:51:51 2002

LSC Report - Integrated Chromatogram

File : C:\HPCHEM\1\DATA\NOV2601\LBA.D
Operator : 209 GALOS
Acquired : 26 Nov 2001 6:31 pm using AcqMethod NP103001
Instrument : GC/MS 209
Sample Name: 11/20 lab blk scan
Misc Info :
Vial Number: 9
Quant File :GCMS1126.RES (RTE Integrator)



LBA.D GCMS1126.M Thu Jan 03 12:51:57 2002

Library Search Compound Report

Data File : C:\HPCHEM\1\DATA\NOV2601\LBA.D
Acq On : 26 Nov 2001 6:31 pm
Sample : 11/20 lab blk scan
Misc :
MS Integration Params: LSCINT.P

Vial: 9
Operator: 209 GALOS
Inst : GC/MS 209
Multiplr: 1.00

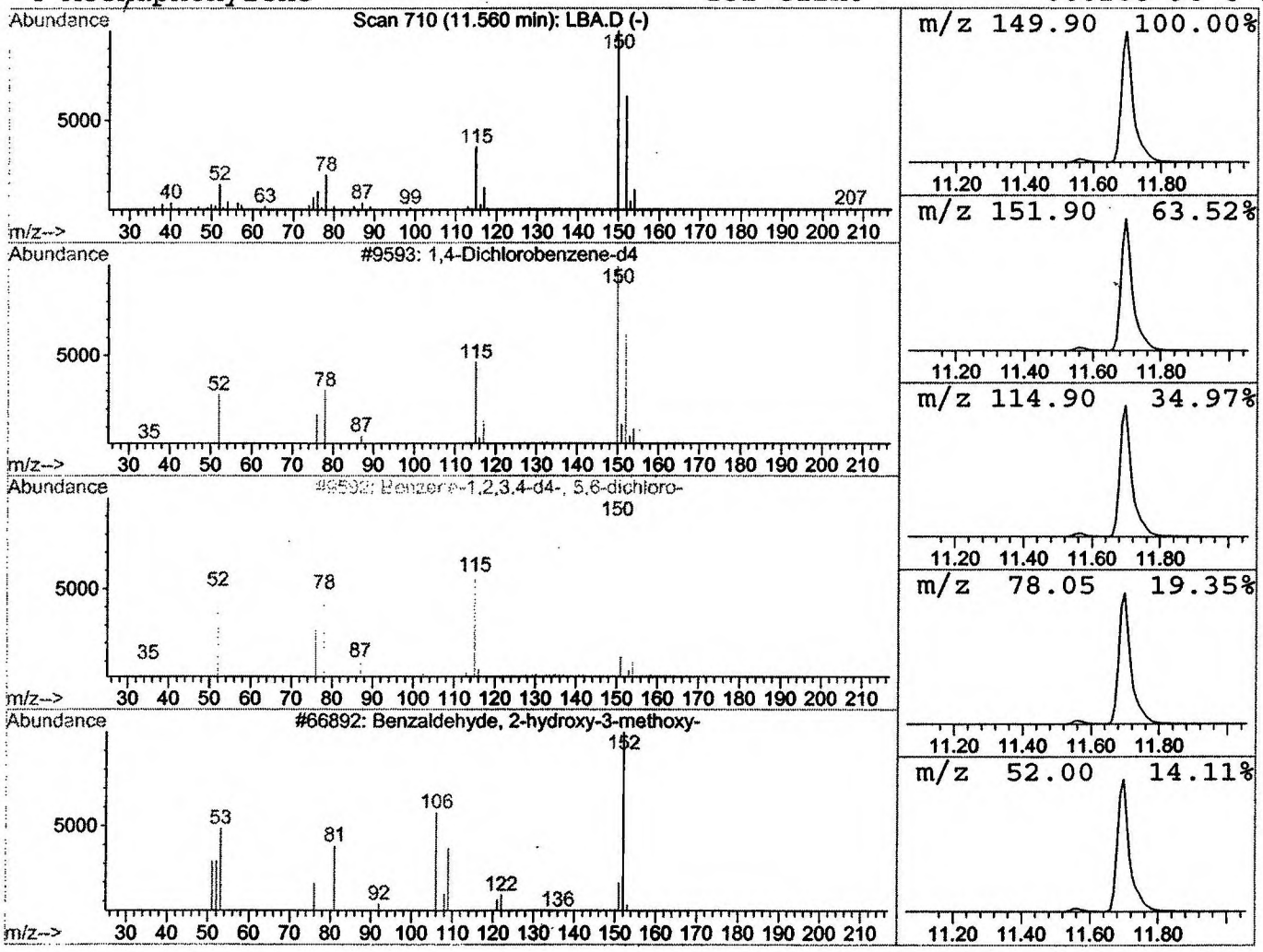
Quant Method : C:\HPCHEM\1\METHODS\GCMS1126.M (RTE Integrator)
Title : 209 625/TCLP calibration table
Library : C:\DATABASE\NBS75K.L

****

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.56	0.51 ug/l	129074	1,4-Dichlorobenzene-d4	11.70

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	25
3	Benzaldehyde, 2-hydroxy-3-methoxy-	152	C8H8O3	000148-53-8	10
4	Acenaphthylene	152	C12H8	000208-96-8	10



Library Search Compound Report

Data File : C:\HPCHEM\1\DATA\NOV2601\LBA.D
Acq On : 26 Nov 2001 6:31 pm
Sample : 11/20 lab blk scan
Misc :
MS Integration Params: LSCINT.P

Vial: 9
Operator: 209 GALOS
Inst : GC/MS 209
Multiplr: 1.00

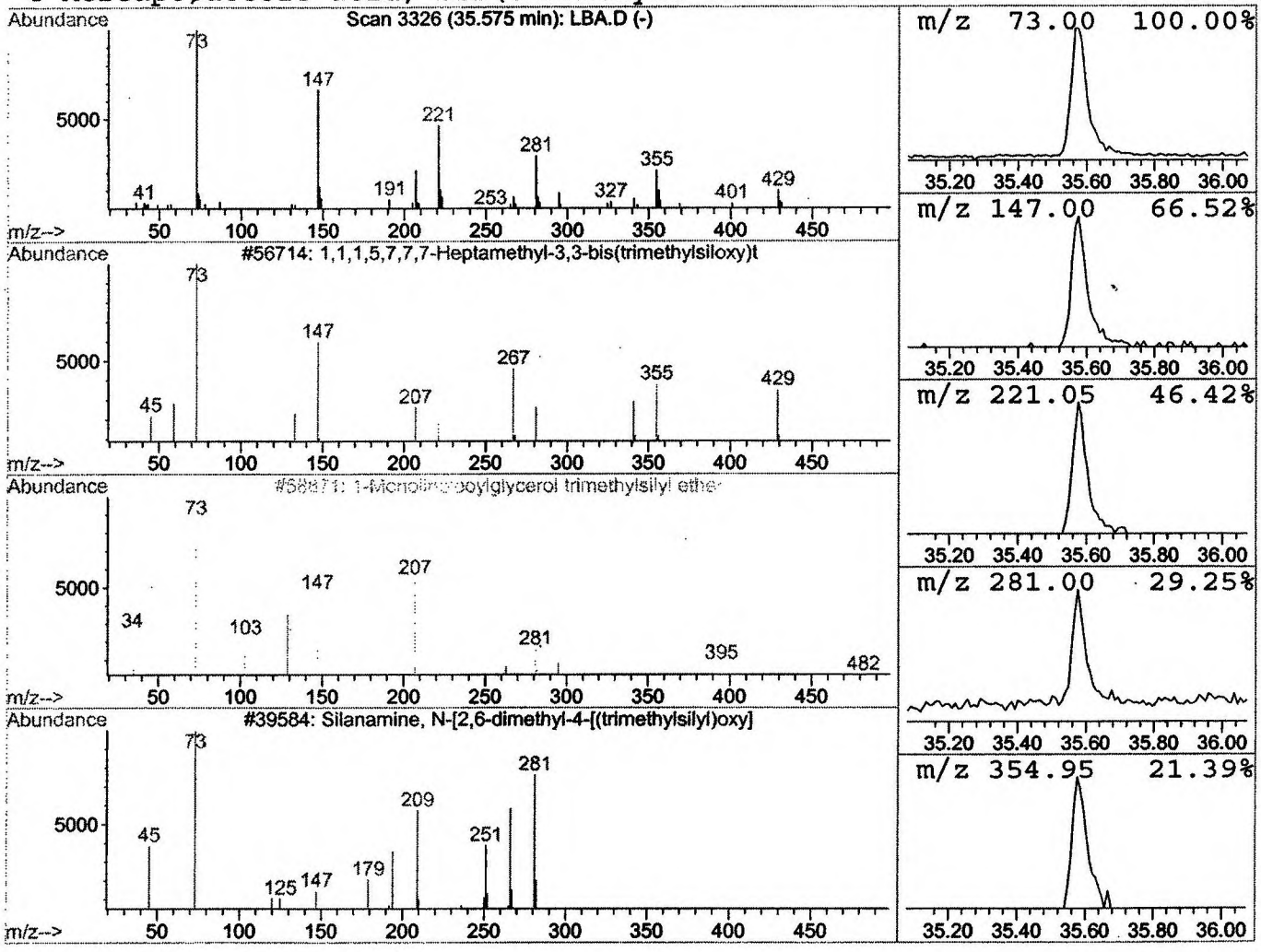
Quant Method : C:\HPCHEM\1\METHODS\GCMS1126.M (RTE Integrator)
Title : 209 625/TCLP calibration table
Library : C:\DATABASE\NBS75K.L

****

Peak Number 2 1,1,1,5,7,7,7-Heptamethyl-3,3- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
35.58	0.59 ug/l	115211	Perylene-d12	37.12

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	32
2	1-Monolinoleoylglycerol trimethylsi	498	C27H54O4Si2	054284-45-6	20
3	Silanamine, N-[2,6-dimethyl-4-[(tri	281	C14H27NOSi2	072088-09-6	11
4	Mercaptoacetic acid, bis(trimethyls	236	C8H20O2SSi2	006398-62-5	10



Library Search Compound Report

Data File : C:\HPCHEM\1\DATA\NOV2601\LBA.D
Acq On : 26 Nov 2001 6:31 pm
Sample : 11/20 lab blk scan
Misc :
MS Integration Params: LSCINT.P

Vial: 9
Operator: 209 GALOS
Inst : GC/MS 209
Multiplr: 1.00

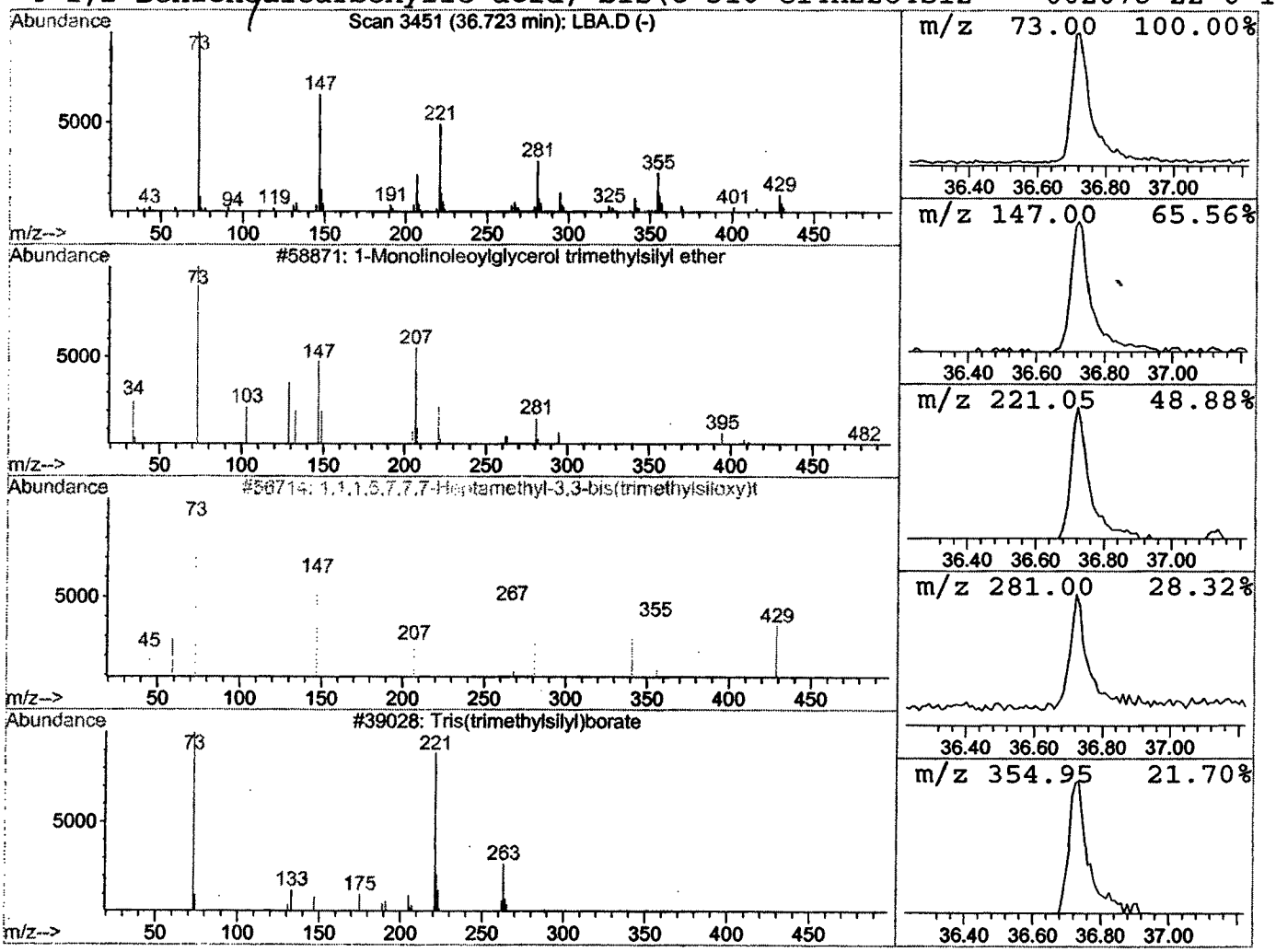
Quant Method : C:\HPCHEM\1\METHODS\GCMS1126.M (RTE Integrator)
Title : 209 625/TCLP calibration table
Library : C:\DATABASE\NBS75K.L

****

Peak Number 3 1-Monolinoleoylglycerol trimet Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
36.72	0.91 ug/l	176851	Perylene-d12	37.12

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Monolinoleoylglycerol trimethylsi	498	C27H54O4Si2	054284-45-6	23
2	1,1,1,5,7,7,7-Heptamethyl-3,3-bis(t	444	C13H40O5Si6	038147-00-1	17
3	Tris(trimethylsilyl)borate	278	C9H27BO3Si3	004325-85-3	12
4	1,2-Benzenedicarboxylic acid, bis(t	310	C14H22O4Si2	002078-22-0	10



Library Search Compound Report

Data File : C:\HPCHEM\1\DATA\NOV2601\LBA.D
Acq On : 26 Nov 2001 6:31 pm
Sample : 11/20 lab blk scan
Misc :
MS Integration Params: LSCINT.P

Vial: 9
Operator: 209 GALOS
Inst : GC/MS 209
Multiplr: 1.00

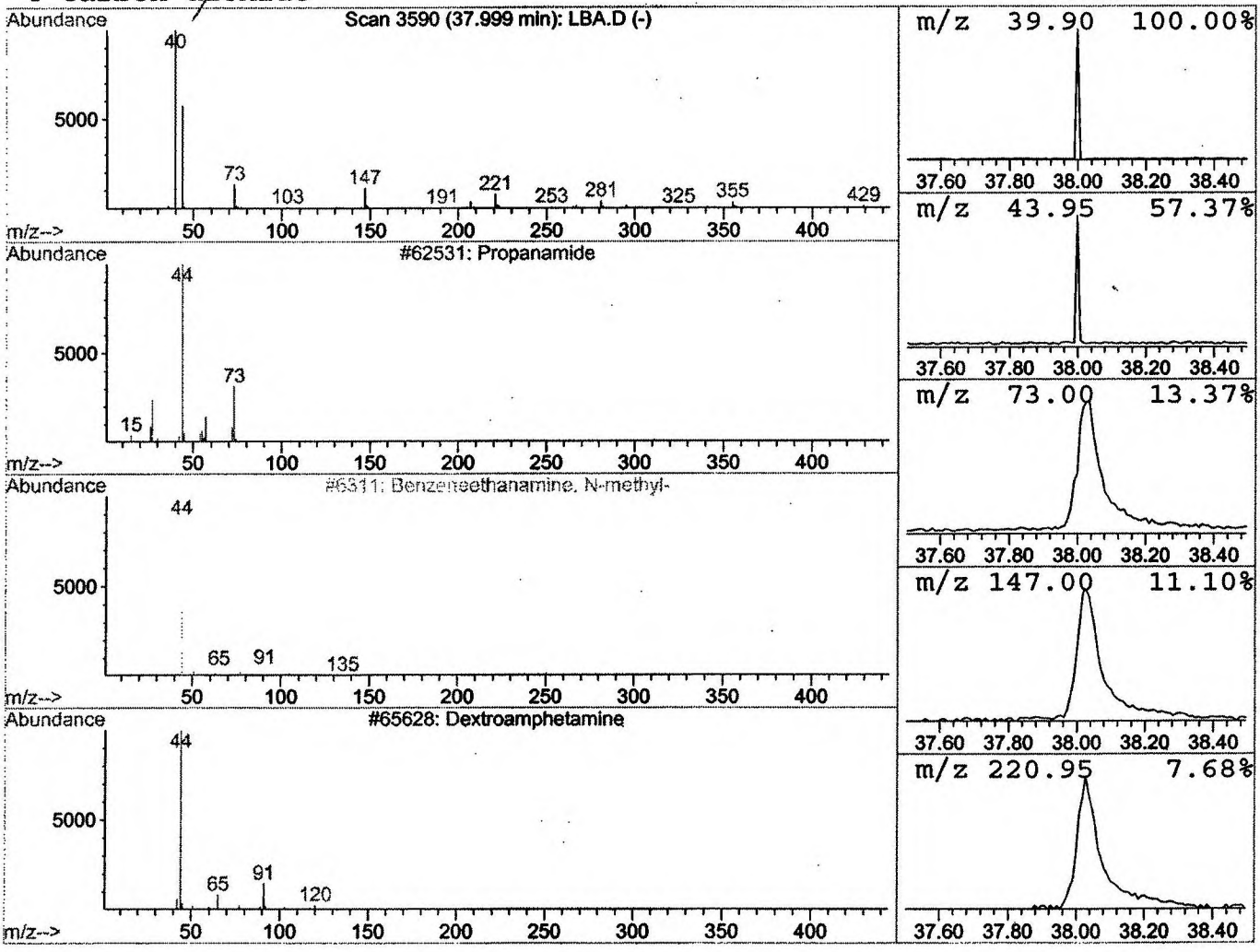
Quant Method : C:\HPCHEM\1\METHODS\GCMS1126.M (RTE Integrator)
Title : 209 625/TCLP calibration table
Library : C:\DATABASE\NBS75K.L

****

Peak Number 4 Propanamide Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
38.00	0.44 ug/l	85848	Perylene-d12	37.12

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Propanamide	73	C3H7NO	000079-05-0	5
2	Benzeneethanamine, N-methyl-	135	C9H13N	000589-08-2	4
3	Dextroamphetamine	135	C9H13N	000051-64-9	3
4	Carbon dioxide	44	CO2	000124-38-9	3



Library Search Compound Report

Data File : C:\HPCHEM\1\DATA\NOV2601\LBA.D
Acq On : 26 Nov 2001 6:31 pm
Sample : 11/20 lab blk scan
Misc :
MS Integration Params: LSCINT.P

Vial: 9
Operator: 209 GALOS
Inst : GC/MS 209
Multiplr: 1.00

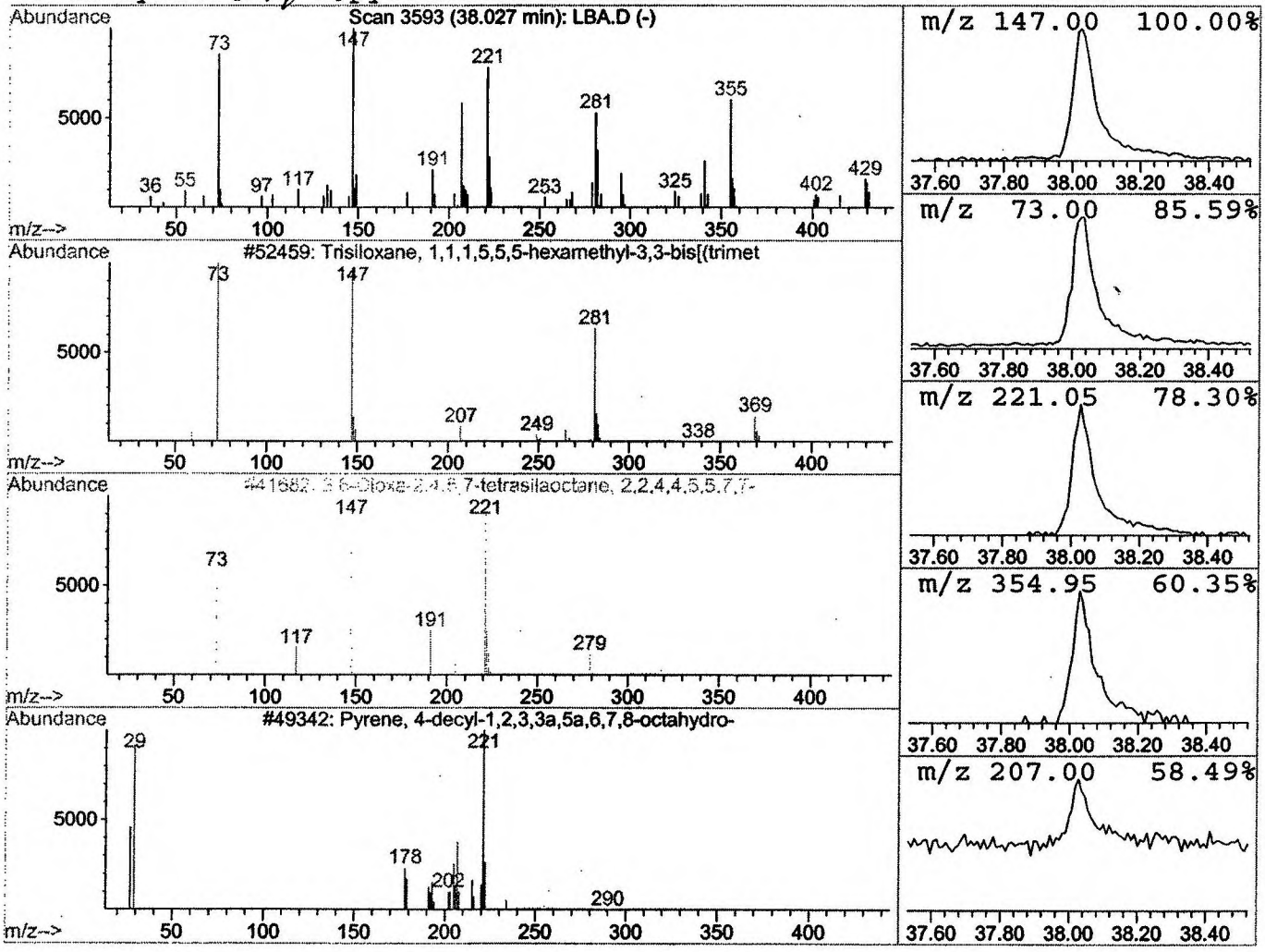
Quant Method : C:\HPCHEM\1\METHODS\GCMS1126.M (RTE Integrator)
Title : 209 625/TCLP calibration table
Library : C:\DATABASE\NBS75K.L

****

Peak Number 5 Trisiloxane, 1,1,1,5,5,5-hexam Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
38.03	1.13 ug/l	219627	Perylene-d12	37.12

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Trisiloxane, 1,1,1,5,5,5-hexamethyl	384	C12H36O4Si5	003555-47-3	12
2	3,6-Dioxa-2,4,5,7-tetrasilaoctane,	294	C10H30O2Si4	004342-25-0	12
3	Pyrene, 4-decyl-1,2,3,3a,5a,6,7,8-o	350	C26H38	055493-75-9	10
4	4H-Pyrido[1,2-a]pyrimidine-3-carbox	221	C11H15N3O2	064399-29-7	10



Library Search Compound Report

Data File : C:\HPCHEM\1\DATA\NOV2601\LBA.D
Acq On : 26 Nov 2001 6:31 pm
Sample : 11/20 lab blk scan
Misc :
MS Integration Params: LSCINT.P

Vial: 9
Operator: 209 GALOS
Inst : GC/MS 209
Multiplr: 1.00

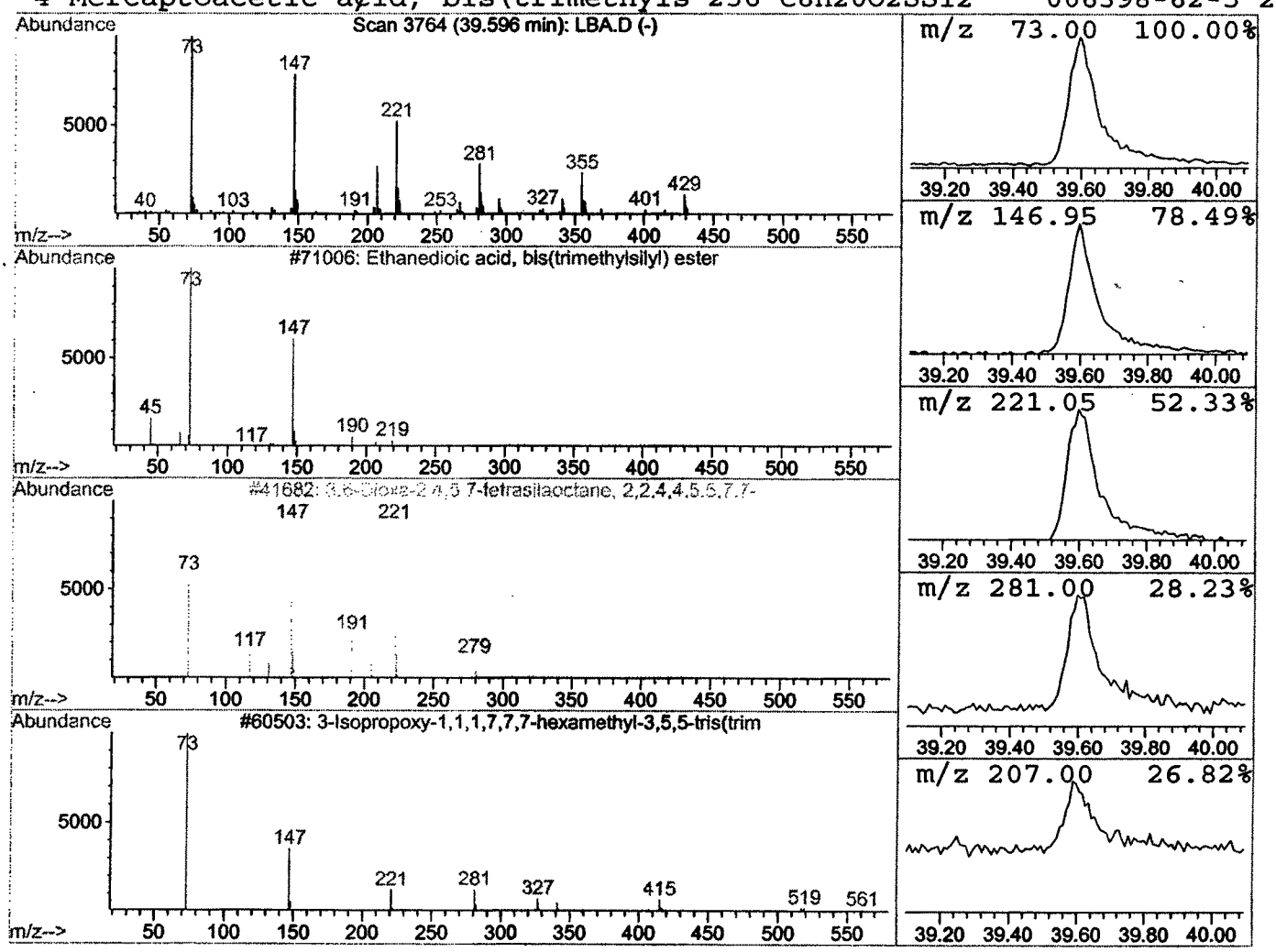
Quant Method : C:\HPCHEM\1\METHODS\GCMS1126.M (RTE Integrator)
Title : 209 625/TCLP calibration table
Library : C:\DATABASE\NBS75K.L

****

Peak Number 6 Ethanedioic acid, bis(trimethy Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
39.60	2.07 ug/l	401611	Perylene-d12	37.12

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Ethanedioic acid, bis(trimethylsily	234	C8H18O4Si2	018294-04-7	27
2	3,6-Dioxa-2,4,5,7-tetrasilaoctane,	294	C10H30O2Si4	004342-25-0	25
3	3-Isopropoxy-1,1,1,7,7,7-hexamethyl	576	C18H52O7Si7	071579-69-6	25
4	Mercaptoacetic acid, bis(trimethyls	236	C8H20O2SSi2	006398-62-5	23



Library Search Compound Report

Data File : C:\HPCHEM\1\DATA\NOV2601\LBA.D
Acq On : 26 Nov 2001 6:31 pm
Sample : 11/20 lab blk scan
Misc :
MS Integration Params: LSCINT.P

Vial: 9
Operator: 209 GALOS
Inst : GC/MS 209
Multiplr: 1.00

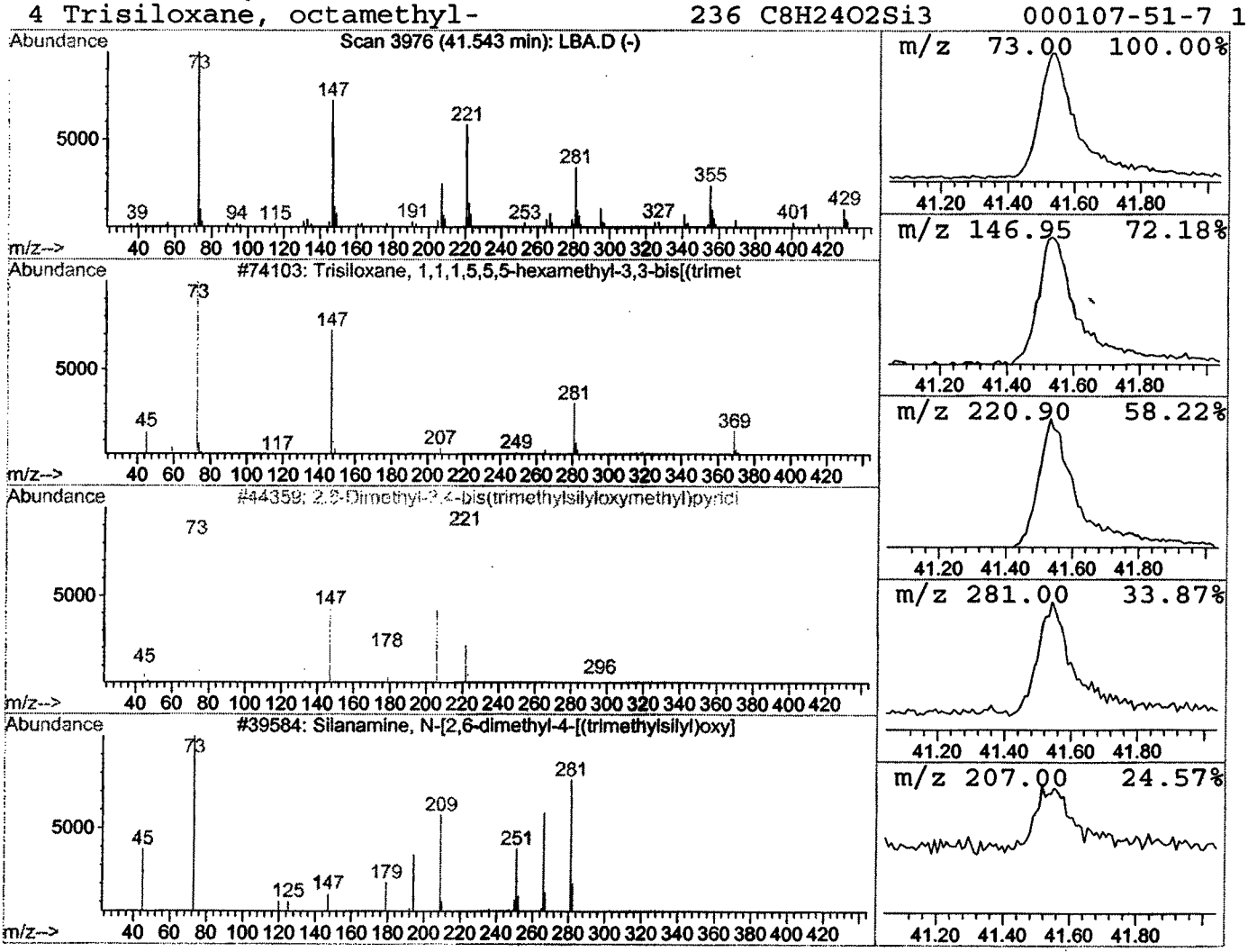
Quant Method : C:\HPCHEM\1\METHODS\GCMS1126.M (RTE Integrator)
Title : 209 625/TCLP calibration table
Library : C:\DATABASE\NBS75K.L

****

Peak Number 7 Trisiloxane, 1,1,1,5,5,5-hexam Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
41.54	1.98 ug/l	383337	Perylene-d12	37.12

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Trisiloxane, 1,1,1,5,5,5-hexamethyl	384	C12H36O4Si5	003555-47-3	30
2	2,6-Dimethyl-3,4-bis(trimethylsilyl	311	C15H29NO2Si2	000000-00-0	12
3	Silanamine, N-[2,6-dimethyl-4-[(tri	281	C14H27NOSi2	072088-09-6	11
4	Trisiloxane, octamethyl-	236	C8H24O2Si3	000107-51-7	10



Library Search Compound Report

Data File : C:\HPCHEM\1\DATA\NOV2601\LBA.D
Acq On : 26 Nov 2001 6:31 pm
Sample : 11/20 lab blk scan
Misc :
MS Integration Params: LSCINT.P

Vial: 9
Operator: 209 GALOS
Inst : GC/MS 209
Multiplr: 1.00

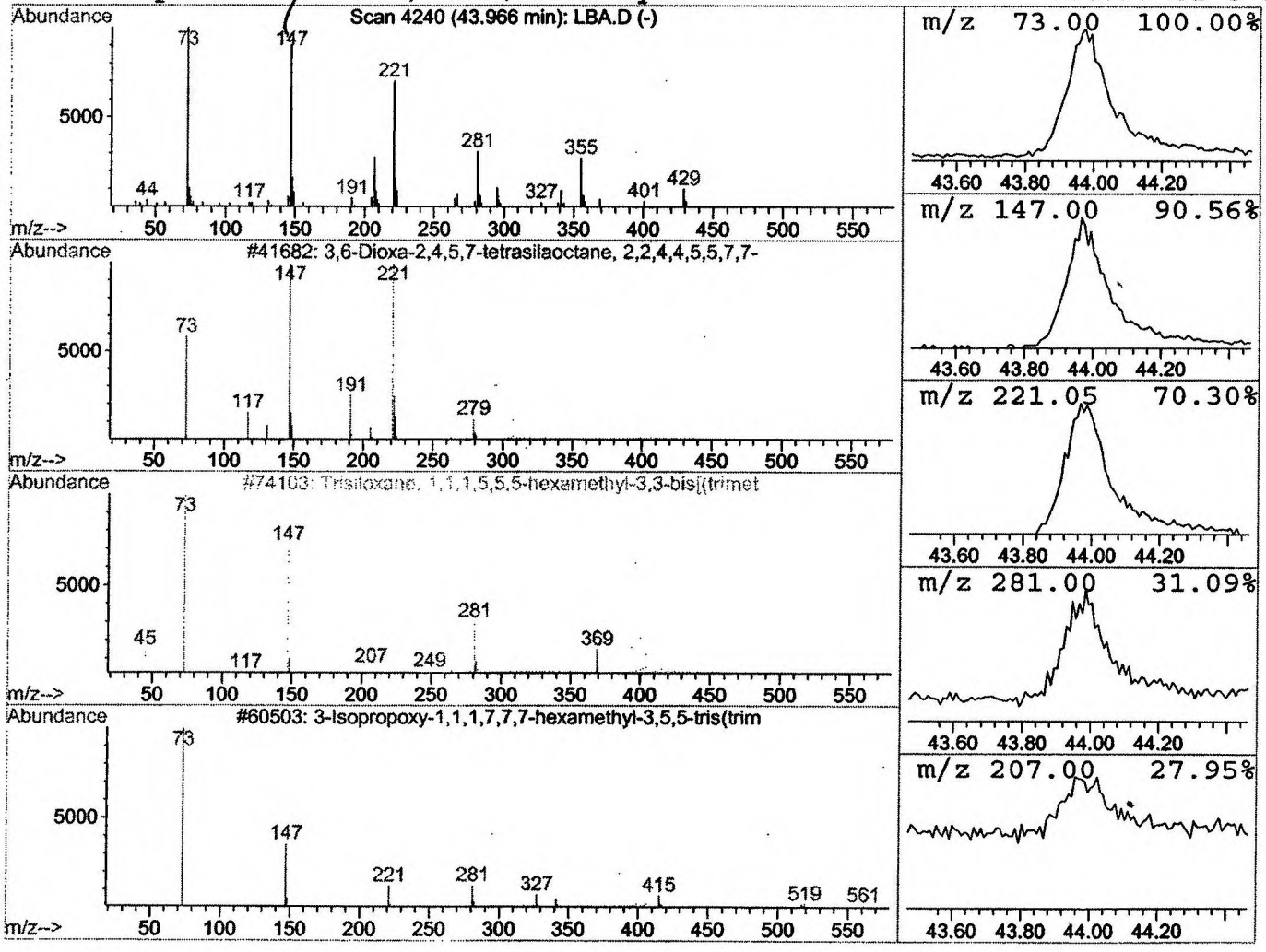
Quant Method : C:\HPCHEM\1\METHODS\GCMS1126.M (RTE Integrator)
Title : 209 625/TCLP calibration table
Library : C:\DATABASE\NBS75K.L

****

Peak Number 8 3,6-Dioxa-2,4,5,7-tetrasilaoct Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
43.97	1.89 ug/l	365657	Perylene-d12	37.12

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, 1,1,1,5,5,5-hexamethyl	384	C12H36O4Si5	003555-47-3	35
3	3-Isopropoxy-1,1,1,7,7,7-hexamethyl	576	C18H52O7Si7	071579-69-6	28
4	Mercaptoacetic acid, bis(trimethyls	236	C8H20O2SSi2	006398-62-5	17



Library Search Compound Report

Data File : C:\HPCHEM\1\DATA\NOV2601\LBA.D
Acq On : 26 Nov 2001 6:31 pm
Sample : 11/20 lab blk scan
Misc :
MS Integration Params: LSCINT.P

Vial: 9
Operator: 209 GALOS
Inst : GC/MS 209
Multiplr: 1.00

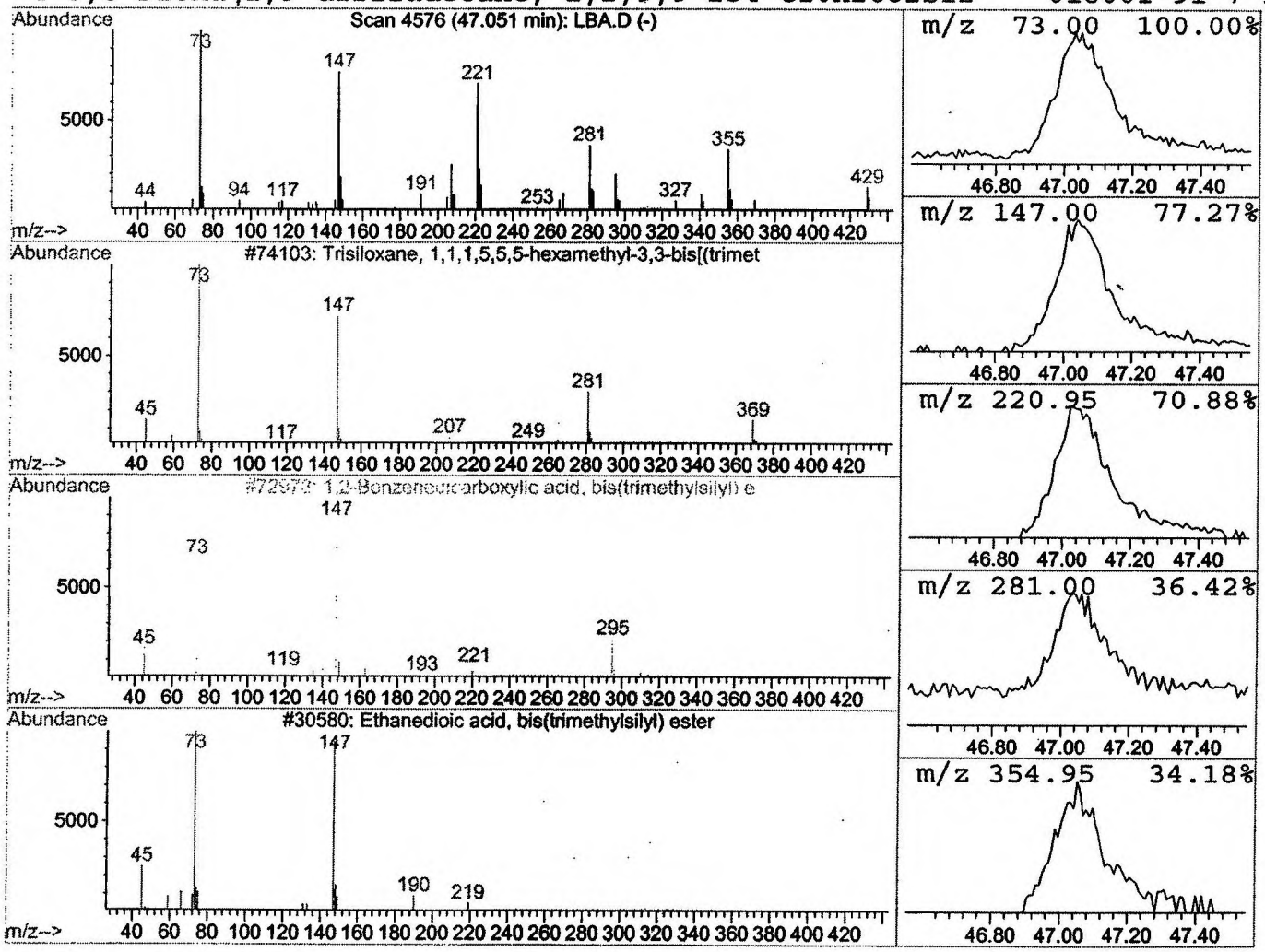
Quant Method : C:\HPCHEM\1\METHODS\GCMS1126.M (RTE Integrator)
Title : 209 625/TCLP calibration table
Library : C:\DATABASE\NBS75K.L

****

Peak Number 9 Trisiloxane, 1,1,1,5,5,5-hexam Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
47.05	0.69 ug/l	133756	Perylene-d12	37.12

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Trisiloxane, 1,1,1,5,5,5-hexamethyl	384	C12H36O4Si5	003555-47-3	35
2	1,2-Benzenedicarboxylic acid, bis(t	310	C14H22O4Si2	002078-22-0	35
3	Ethanedioic acid, bis(trimethylsily	234	C8H18O4Si2	018294-04-7	12
4	3,8-Dioxa-2,9-disiladecane, 2,2,9,9	234	C10H26O2Si2	018001-91-7	12



Tentatively Identified Compound (LSC) summary

Operator ID: 209 GALOS Date Acquired: 26 Nov 2001 6:31 pm
Data File: C:\HPCHEM\1\DATA\NOV2601\LBA.D
Name: 11/20 lab blk scan
Misc:
Method: C:\HPCHEM\1\METHODS\GCMS1126.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
1,4-Dichlorobenzene-	11.56	0.5	ug/l	129074	ISTD01	11.70	5101610	20.0
1,1,1,5,7,7,7-Heptam	35.58	0.6	ug/l	115211	ISTD06	37.12	3876040	20.0
1-Monolinoleoylglyce	36.72	0.9	ug/l	176851	ISTD06	37.12	3876040	20.0
Propanamide	38.00	0.4	ug/l	85848	ISTD06	37.12	3876040	20.0
Trisiloxane, 1,1,1,5	38.03	1.1	ug/l	219627	ISTD06	37.12	3876040	20.0
Ethanedioic acid, bi	39.60	2.1	ug/l	401611	ISTD06	37.12	3876040	20.0
Trisiloxane, 1,1,1,5	41.54	2.0	ug/l	383337	ISTD06	37.12	3876040	20.0
3,6-Dioxa-2,4,5,7-te	43.97	1.9	ug/l	365657	ISTD06	37.12	3876040	20.0
Trisiloxane, 1,1,1,5	47.05	0.7	ug/l	133756	ISTD06	37.12	3876040	20.0

LBA.D GCMS1126.M Thu Jan 03 12:52:42 2002

Column Fleet