User: Galos, Marie
Westchester Dept. of Labs & Research
Date: 04/18/2002 Time: 12:35:01

Sample: AE06430
Analysis: $B1GCMS Blank for GCMS Scan
Units: ug
Started: 4/17/20 00:09 Ended: 4/17/20 00:09 Analyst: MG
Result source: manual entry
Page: 1

	Component Name	Viol	Result	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	2-Chloronaphthalene		< MDL	2.0
16	Dimethylphthalate		< MDL	2.0
17	2,6-Dinitrotoluene		< MDL	2.0
18	Acenaphthylene		< MDL	2.0
19	Acenaphthene		< MDL	2.0
20	2,4-Dinitrotoluene		< MDL	2.0
21	Diethylphthalate		< MDL	2.0
22	4-Chlorophenylphenylether		< MDL	2.0
23	Fluorene		< MDL	2.0
24	Azobenzene		< MDL	2.0
25	n-Nitrosodiphenylamine		< MDL	2.0
26	4-Bromophenylphenylether		< MDL	2.0
27	Hexachlorobenzene		< MDL	2.0
28	Phenanthrene		< MDL	2.0
29	Anthracene		< MDL	2.0
30	Di-n-butylphthalate		< MDL	2.0
31	Fluoranthene		< MDL	2.0
32	Pyrene		< MDL	2.0
33	Butylbenzylphthalate		< MDL	2.0
34	Benzo(a)anthracene		< MDL	2.0
35	bis(2-Ethylhexyl)phthalate		< MDL	2.0
36	Chrysene		< MDL	2.0
37	Di-n-octylphthalate		< MDL	2.0
38	Benzo(b)fluoranthene		< MDL	2.0
39	Benzo(k)fluoranthene		< MDL	2.0
40	Benzo(a)pyrene		< MDL	2.0
41	Indeno(1,2,3-cd)pyrene		< MDL	2.0
42	Dibenz(a,h)anthracene		< MDL	2.0
43	Benzo(g,h,i)perylene		< MDL	2.0

Quantitation Report (QT Reviewed)

Data File : C:\HPCHEM\1\DATA\APR1702\LB.D Vial: 3
Acq On : 17 Apr 2002 10:13 am Operator:
Sample : gc/ms scan 3/19 Inst : 209
Misc : Multiplr: 1.00
MS Integration Params: GOOFY.P
Quant Time: Apr 18 10:55 2002 Quant Results File: GCMS0412.RES

Quant Method : C:\HPCHEM\1\METHODS\GCMS0412.M (RTE Integrator)
Title : 209 625/TCLP calibration table
Last Update : Wed Apr 17 11:48:30 2002
Response via : Initial Calibration
DataAcq Meth : GCMS0412

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	12.21	152	502246	20.00	ug/l	-0.03
10) Naphthalene-d8	15.84	136	1798780	20.00	ug/l	-0.03
17) Acenaphthene-d10	21.15	164	1002923	20.00	ug/l	-0.02
30) Phenanthrene-d10	25.58	188	1373846	20.00	ug/l	-0.03
39) Chrysene-d12	33.64	240	800977	20.00	ug/l	-0.04
45) Perylene-d12	37.87	264	534659	20.00	ug/l	-0.05
System Monitoring Compounds						
28) TCMX	0.00	209	0	0.00	ug/L	
Spiked Amount 10.000			Recovery = 0.00%			
38) 2,4-DDD	30.66	235	100629	6.29	ug/mL	0.16
Spiked Amount 5.000			Recovery = 125.80%			
Target Compounds						Qvalue
2) N-nitroso-dimethyl amine	5.46	74	286	N.D.		
3) bis(2-Chloroethyl)ether	0.00	93	0	N.D.		
4) 1,3-Dichlorobenzene	12.12	146	541	N.D.		
5) 1,4-Dichlorobenzene	12.25	146	521	N.D.		
6) 1,2-Dichlorobenzene	12.77	146	105	N.D.		
7) 2,2'-oxybis(1-Chloropropan	0.00	45	0	N.D.		
8) N-Nitroso-di-n-propylamine	0.00	70	0	N.D.		
9) Hexachloroethane	0.00	117	0	N.D.		
11) Nitrobenzene	13.43	77	471	N.D.		
12) Isophorone	0.00	82	0	N.D.		
13) bis(2-Chloroethoxy)methane	0.00	93	0	N.D.		
14) 1,2,4-Trichlorobenzene	0.00	180	0	N.D.		
15) Naphthalene	15.90	128	714	N.D.		
16) Hexachlorobutadiene	16.44	225	903	N.D.		
18) Hexachlorocyclopentadiene	0.00	237	0	N.D.		
19) 2-Chloronaphthalene	0.00	162	0	N.D.		
20) Dimethylphthalate	0.00	163	0	N.D.		
21) 2,6-Dinitrotoluene	0.00	165	0	N.D.		
22) Acenaphthylene	0.00	152	0	N.D.		
23) Acenaphthene	0.00	153	0	N.D.		
24) 2,4-Dinitrotoluene	0.00	165	0	N.D.		
25) Diethylphthalate	22.63	149	1396	N.D.		
26) 4-Chlorophenyl-phenylether	0.00	204	0	N.D.		
27) Fluorene	0.00	166	0	N.D.		
29) Azobenzen/ 1,2-Diphenylhyd	0.00	77	0	N.D.		
31) N-Nitrosodiphenylamine	0.00	168	0	N.D.		
32) 4-Bromophenyl-phenylether	0.00	248	0	N.D.		

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

Quantitation Report (QT Reviewed)

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Acq On : 17 Apr 2002 10:13 am Operator:
Sample : gc/ms scan 3/19 Inst : 209
Misc : Multiplr: 1.00
MS Integration Params: GOOFY.P
Quant Time: Apr 18 10:55 2002 Quant Results File: GCMS0412.RES

Quant Method : C:\HPCHEM\1\METHODS\GCMS0412.M (RTE Integrator)
Title : 209 625/TCLP calibration table
Last Update : Wed Apr 17 11:48:30 2002
Response via : Initial Calibration
DataAcq Meth : GCMS0412

	Compound	R.T.	QIon	Response	Conc Unit	Qvalue
33)	Hexachlorobenzene	0.00	284	0	N.D.	
34)	Phenanthrene	25.58	178	391	N.D.	
35)	Anthracene	25.58	178	307	N.D.	
36)	Di-n-butylphthalate	27.56	149	16010	0.36 ug/l	98
37)	Fluoranthene	29.28	202	106	N.D.	
40)	Pyrene	29.95	202	372	N.D.	
41)	Butylbenzylphthalate	32.08	149	1142	N.D.	
42)	Benzo(a)anthracene	33.64	228	3415	N.D.	
43)	Bis(2-ethylhexyl)phthalate	33.86	149	8803	0.44 ug/l #	96
44)	Chrysene	33.71	228	1852	N.D.	
46)	Di-n-Octylphthalate	35.59	149	1613	N.D.	
47)	Benzo(b)fluoranthene	36.68	252	2672	N.D.	
48)	Benzo(k)fluoranthene	36.75	252	3384	N.D.	
49)	Benzo(a)pyrene	37.67	252	1725	N.D.	
50)	Indeno(1,2,3-cd)pyrene	0.00	276	0	N.D. d	
51)	Dibenz(a,h)anthracene	0.00	278	0	N.D. d	
52)	Benzo(g,h,i)perylene	0.00	276	0	N.D. d	

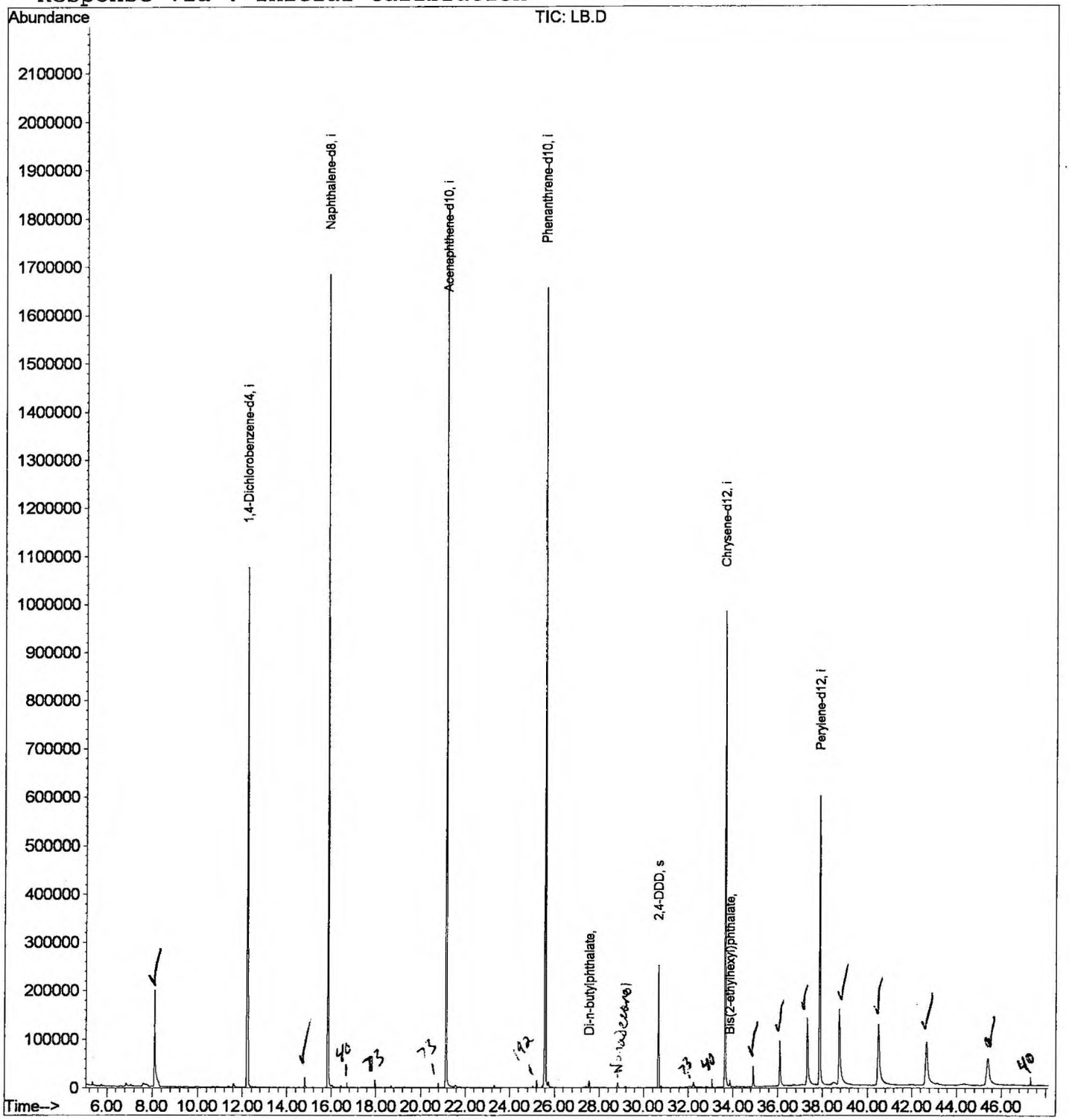
(#) = qualifier out of range (m) = manual integration

Quantitation Report

Data File : C:\HPCHEM\1\DATA\APR1702\LB.D
Acq On : 17 Apr 2002 10:13 am
Sample : gc/ms scan 3/19
Misc :
MS Integration Params: GOOFY.P
Quant Time: Apr 18 10:55 2002

Vial: 3
Operator:
Inst : 209
Multiplr: 1.00
Quant Results File: GCMS0412.RES

Method : C:\HPCHEM\1\METHODS\GCMS0412.M (RTE Integrator)
Title : 209 625/TCLP calibration table
Last Update : Wed Apr 17 11:48:30 2002
Response via : Initial Calibration



LSC Area Percent Report

Data File : C:\HPCHEM\1\DATA\APR1702\LB.D Vial: 3
Acq On : 17 Apr 2002 10:13 am Operator:
Sample : gc/ms scan 3/19 Inst : 209
Misc : Multiplr: 1.00
MS Integration Params: LSCINT.P

Method : C:\HPCHEM\1\METHODS\GCMS0412.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

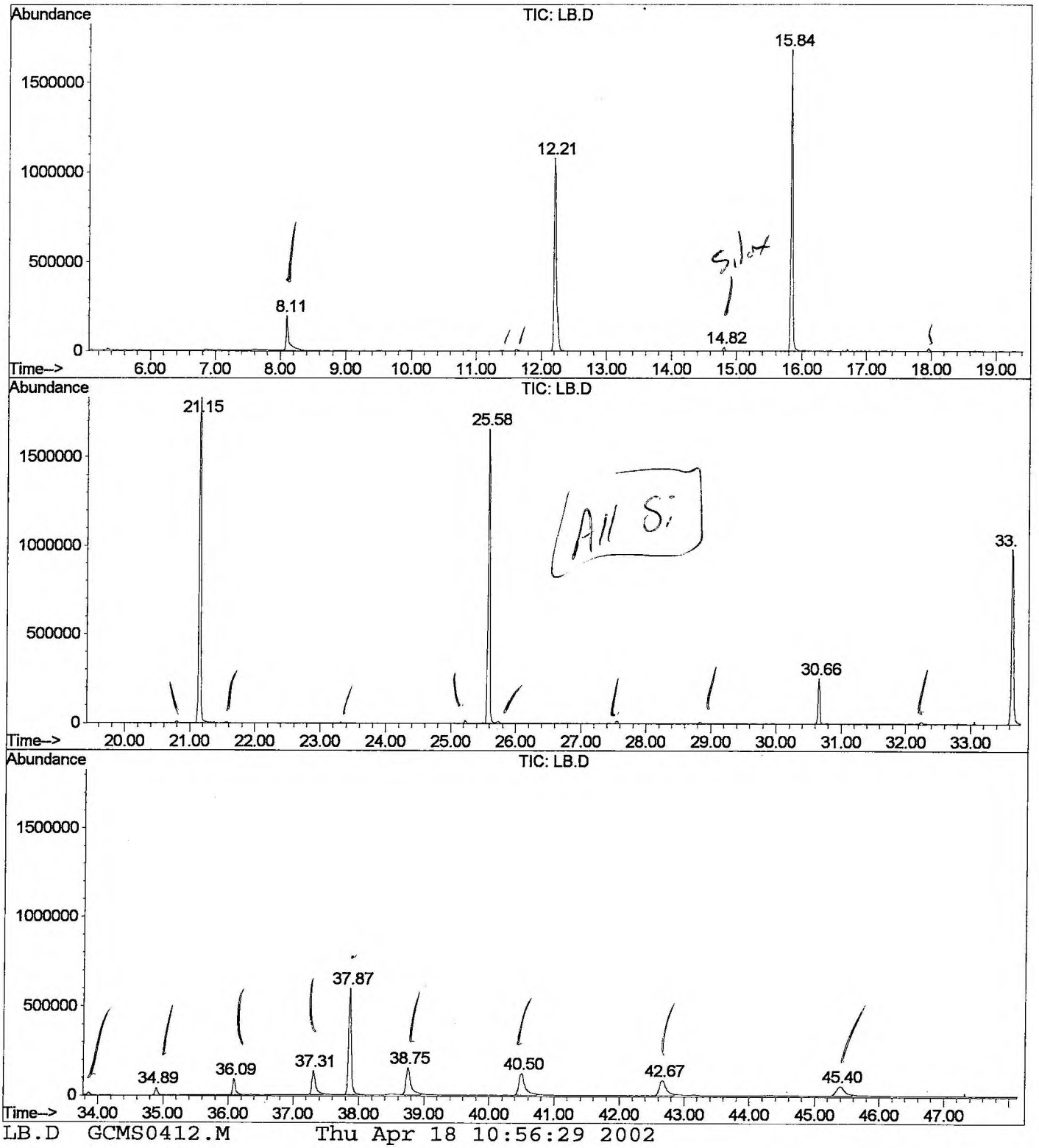
peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	peak area	peak % max.	% of total
1	8.105	329	333	359	rVB	198306	578215	15.34%	2.614%
2	12.209	775	780	797	rVB	1074684	2674292	70.95%	12.091%
3	14.816	1059	1064	1069	rBV2	21673	40517	1.07%	0.183%
4	15.844	1170	1176	1193	rBV	1683676	3372191	89.47%	15.246%
5	21.150	1747	1754	1763	rBV	1828843	3769046	100.00%	17.040%
6	25.585	2230	2237	2248	rBV2	1654776	3419840	90.73%	15.461%
7	30.661	2784	2790	2797	rVB	251098	483151	12.82%	2.184%
8	33.636	3103	3114	3127	rBV2	984008	2335016	61.95%	10.557%
9	34.893	3243	3251	3261	rBV	42356	111628	2.96%	0.505%
10	36.087	3375	3381	3395	rBV	92002	278661	7.39%	1.260%
11	37.308	3507	3514	3531	rBV2	135647	537375	14.26%	2.430%
12	37.868	3567	3575	3597	rBV2	596786	1791450	47.53%	8.099%
13	38.749	3662	3671	3693	rBV	154362	788207	20.91%	3.564%
14	40.503	3850	3862	3888	rBV2	123718	788043	20.91%	3.563%
15	42.669	4083	4098	4118	rBV4	87808	651983	17.30%	2.948%
16	45.396	4377	4395	4415	rBV5	54836	498841	13.24%	2.255%

Sum of corrected areas: 22118456

LB.D GCMS0412.M Thu Apr 18 10:56:28 2002

LSC Report - Integrated Chromatogram

File : C:\HPCHEM\1\DATA\APR1702\LB.D
Operator :
Acquired : 17 Apr 2002 10:13 am using AcqMethod GCMS0412
Instrument : 209
Sample Name: gc/ms scan 3/19
Misc Info :
Vial Number: 3
Quant File :GCMS0412.RES (RTE Integrator)



LB.D GCMS0412.M Thu Apr 18 10:56:29 2002

Library Search Compound Report

Data File : C:\HPCHEM\1\DATA\APR1702\LB.D
Acq On : 17 Apr 2002 10:13 am
Sample : gc/ms scan 3/19
Misc :
MS Integration Params: LSCINT.P

Vial: 3
Operator:
Inst : 209
Multiplr: 1.00

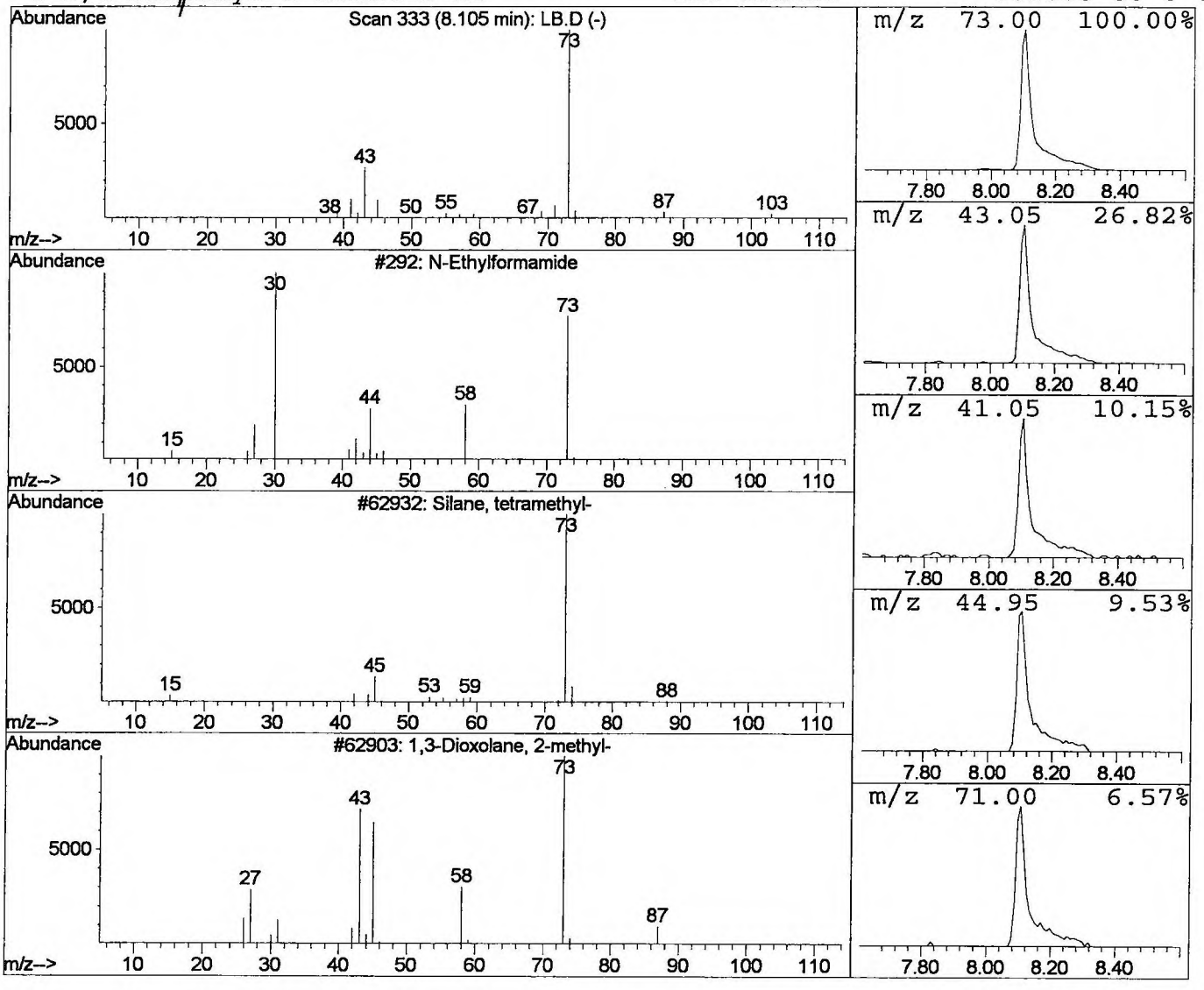
Quant Method : C:\HPCHEM\1\METHODS\GCMS0412.M (RTE Integrator)
Title : 209 625/TCLP calibration table
Library : C:\DATABASE\NBS75K.L

****

Peak Number 1 N-Ethylformamide Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.11	4.32 ug/l	578215	1,4-Dichlorobenzene-d4	12.21

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	N-Ethylformamide	73	C3H7NO	000627-45-2	9
2	Silane, tetramethyl-	88	C4H12Si	000075-76-3	9
3	1,3-Dioxolane, 2-methyl-	88	C4H8O2	000497-26-7	7
4	3,4-Dimethyl-5-hexen-3-ol	128	C8H16O	000000-00-0	4



Library Search Compound Report

Data File : C:\HPCHEM\1\DATA\APR1702\LB.D Vial: 3
Acq On : 17 Apr 2002 10:13 am Operator:
Sample : gc/ms scan 3/19 Inst : 209
Misc : Multiplr: 1.00
MS Integration Params: LSCINT.P

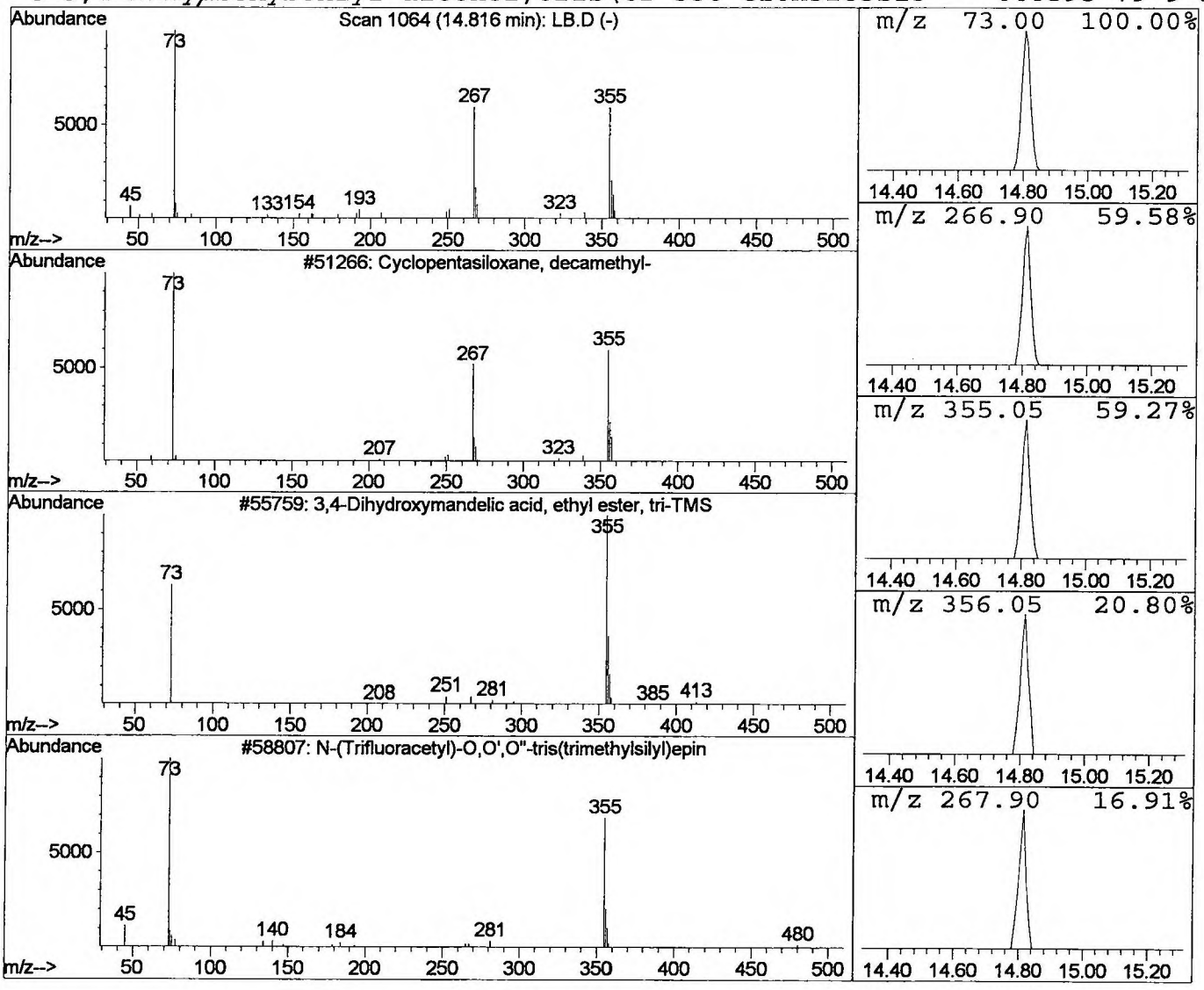
Quant Method : C:\HPCHEM\1\METHODS\GCMS0412.M (RTE Integrator)
Title : 209 625/TCLP calibration table
Library : C:\DATABASE\NBS75K.L

****

Peak Number 2 Cyclopentasiloxane, decamethyl Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.82	0.24 ug/l	40517	Naphthalene-d8	15.84

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclopentasiloxane, decamethyl-	370	C10H30O5Si5	000541-02-6	91
2	3,4-Dihydroxymandelic acid, ethyl e	428	C19H36O5Si3	000000-00-0	40
3	N-(Trifluoracetyl)-O,O',O"-tris(tri	495	C20H36F3NO4Si3	000000-00-0	38
4	3,4-Dihydroxybenzyl alcohol,tris(tr	356	C16H32O3Si3	068595-79-9	38



Library Search Compound Report

Data File : C:\HPCHEM\1\DATA\APR1702\LB.D Vial: 3
Acq On : 17 Apr 2002 10:13 am Operator:
Sample : gc/ms scan 3/19 Inst : 209
Misc : Multiplr: 1.00
MS Integration Params: LSCINT.P

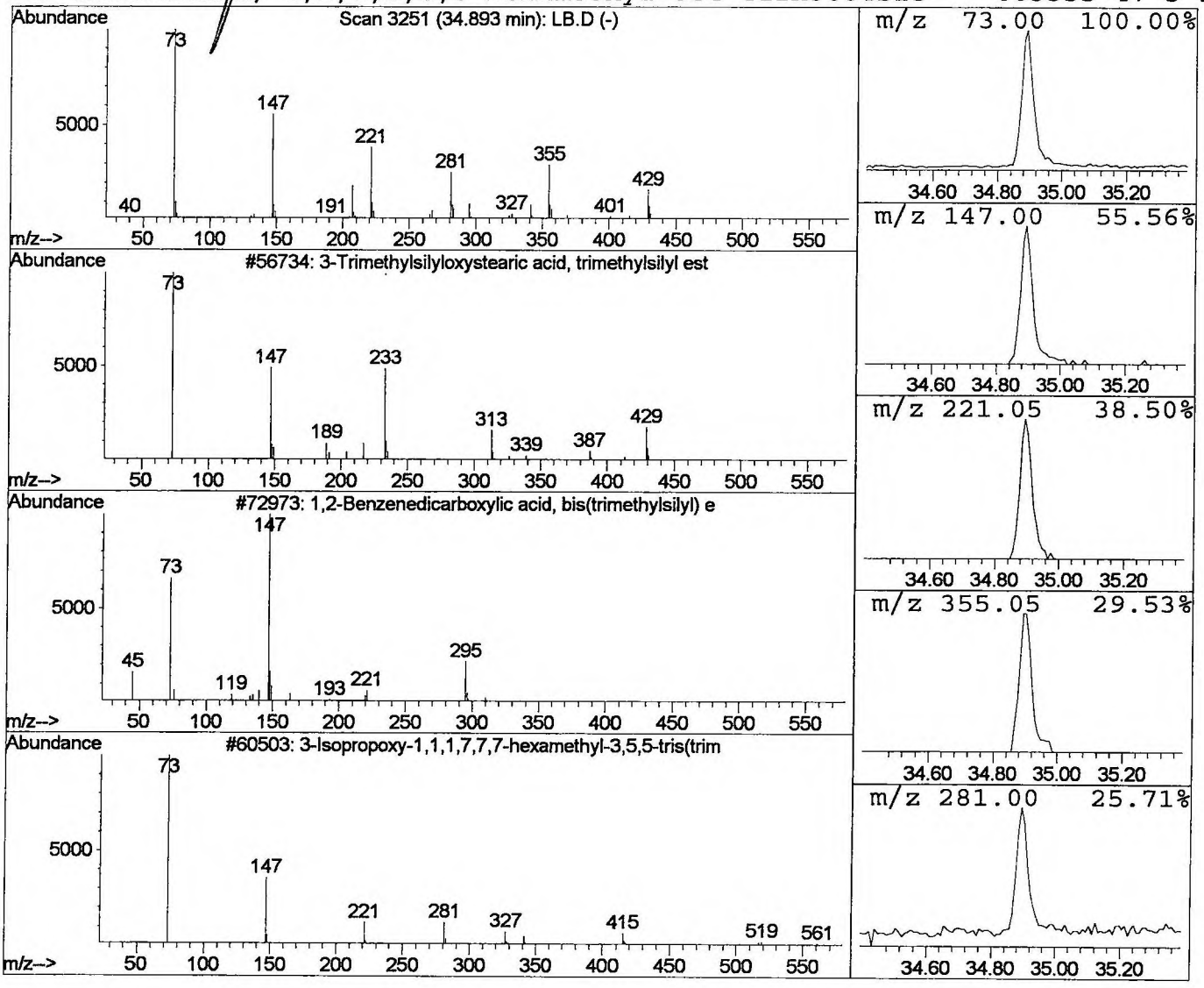
Quant Method : C:\HPCHEM\1\METHODS\GCMS0412.M (RTE Integrator)
Title : 209 625/TCLP calibration table
Library : C:\DATABASE\NBS75K.L

****

Peak Number 3 · 3-Trimethylsilyloxystearic aci Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
34.89	0.96 ug/l	111628	Chrysene-d12	33.64

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	3-Trimethylsilyloxystearic acid, tr	444	C24H52O3Si2	000000-00-0	25
2	1,2-Benzenedicarboxylic acid, bis(t	310	C14H22O4Si2	002078-22-0	16
3	3-Isopropoxy-1,1,1,7,7,7-hexamethyl	576	C18H52O7Si7	071579-69-6	16
4	Trisiloxane, 1,1,1,5,5,5-hexamethyl	384	C12H36O4Si5	003555-47-3	12



Library Search Compound Report

Data File : C:\HPCHEM\1\DATA\APR1702\LB.D Vial: 3
Acq On : 17 Apr 2002 10:13 am Operator:
Sample : gc/ms scan 3/19 Inst : 209
Misc : Multiplr: 1.00
MS Integration Params: LSCINT.P

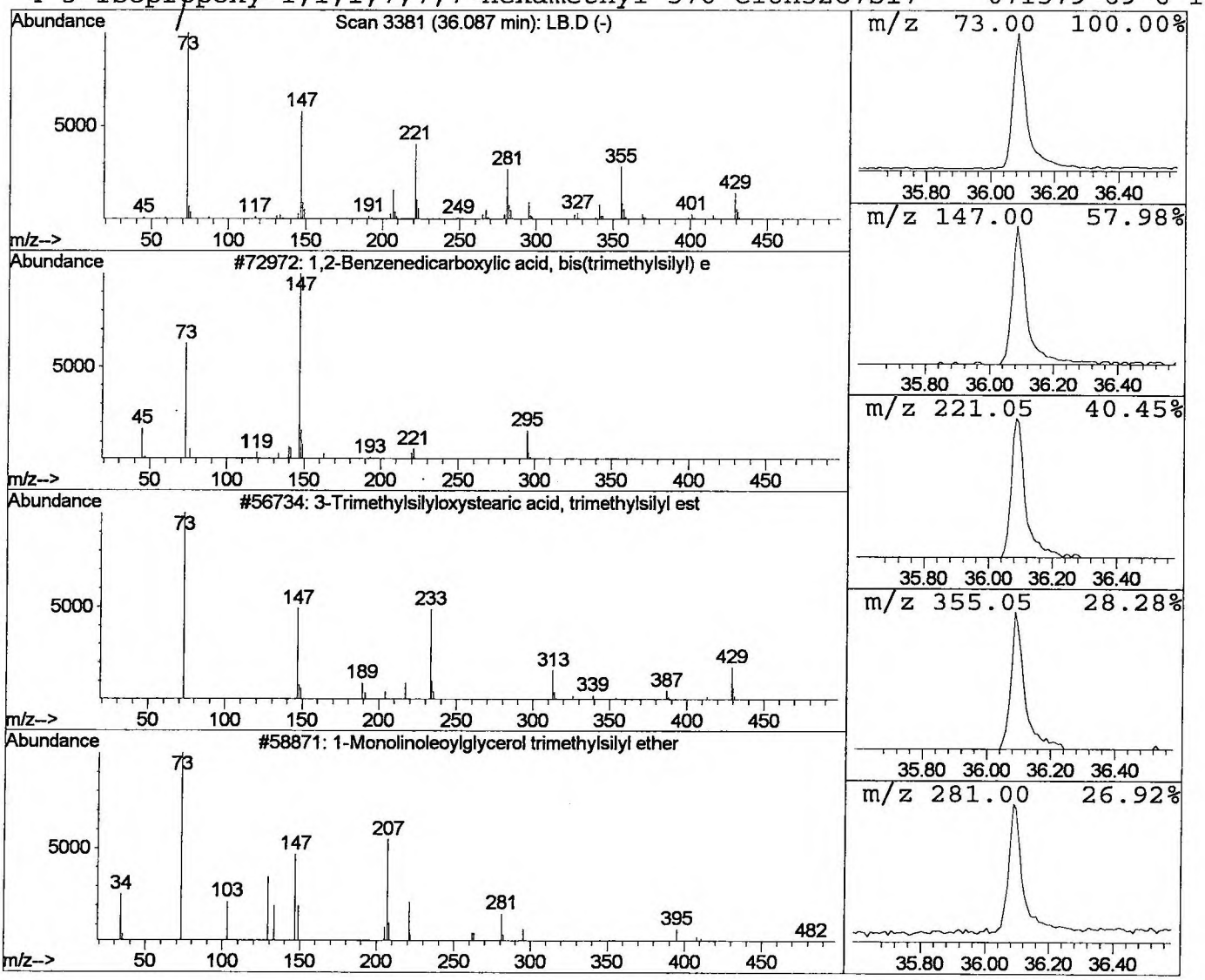
Quant Method : C:\HPCHEM\1\METHODS\GCMS0412.M (RTE Integrator)
Title : 209 625/TCLP calibration table
Library : C:\DATABASE\NBS75K.L

****

Peak Number 4 1,2-Benzenedicarboxylic acid, Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
36.09	3.11 ug/l	278661	Perylene-d12	37.87

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,2-Benzenedicarboxylic acid, bis(t	310	C14H22O4Si2	002078-22-0	27
2	3-Trimethylsilyloxystearic acid, tr	444	C24H52O3Si2	000000-00-0	25
3	1-Monolinoleoylglycerol trimethylsi	498	C27H54O4Si2	054284-45-6	25
4	3-Isopropoxy-1,1,1,7,7,7-hexamethyl	576	C18H52O7Si7	071579-69-6	12



Library Search Compound Report

Data File : C:\HPCHEM\1\DATA\APR1702\LB.D Vial: 3
Acq On : 17 Apr 2002 10:13 am Operator:
Sample : gc/ms scan 3/19 Inst : 209
Misc : Multiplr: 1.00
MS Integration Params: LSCINT.P

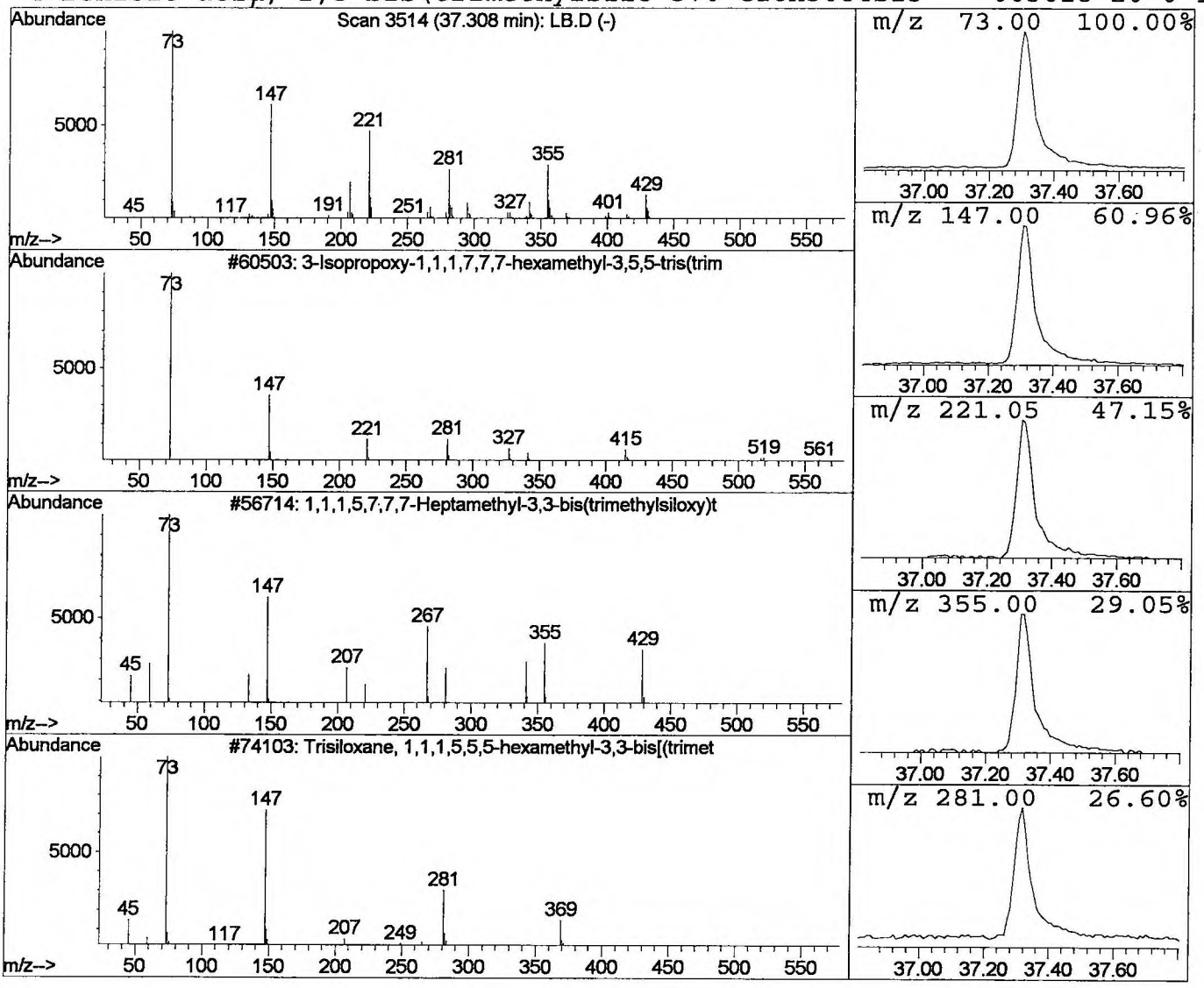
Quant Method : C:\HPCHEM\1\METHODS\GCMS0412.M (RTE Integrator)
Title : 209 625/TCLP calibration table
Library : C:\DATABASE\NBS75K.L

****

Peak Number 5 3-Isopropoxy-1,1,1,7,7,7-hexam Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
37.31	6.00 ug/l	537375	Perylene-d12	37.87

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	3-Isopropoxy-1,1,1,7,7,7-hexamethyl	576	C18H52O7Si7	071579-69-6	32
2	1,1,1,5,7,7,7-Heptamethyl-3,3-bis(t	444	C13H40O5Si6	038147-00-1	28
3	Trisiloxane, 1,1,1,5,5,5-hexamethyl	384	C12H36O4Si5	003555-47-3	12
4	Benzoic acid, 2,5-bis(trimethylsilo	370	C16H30O4Si3	003618-20-0	10



Library Search Compound Report

Data File : C:\HPCHEM\1\DATA\APR1702\LB.D
Acq On : 17 Apr 2002 10:13 am
Sample : gc/ms scan 3/19
Misc :
MS Integration Params: LSCINT.P

Vial: 3
Operator:
Inst : 209
Multiplr: 1.00

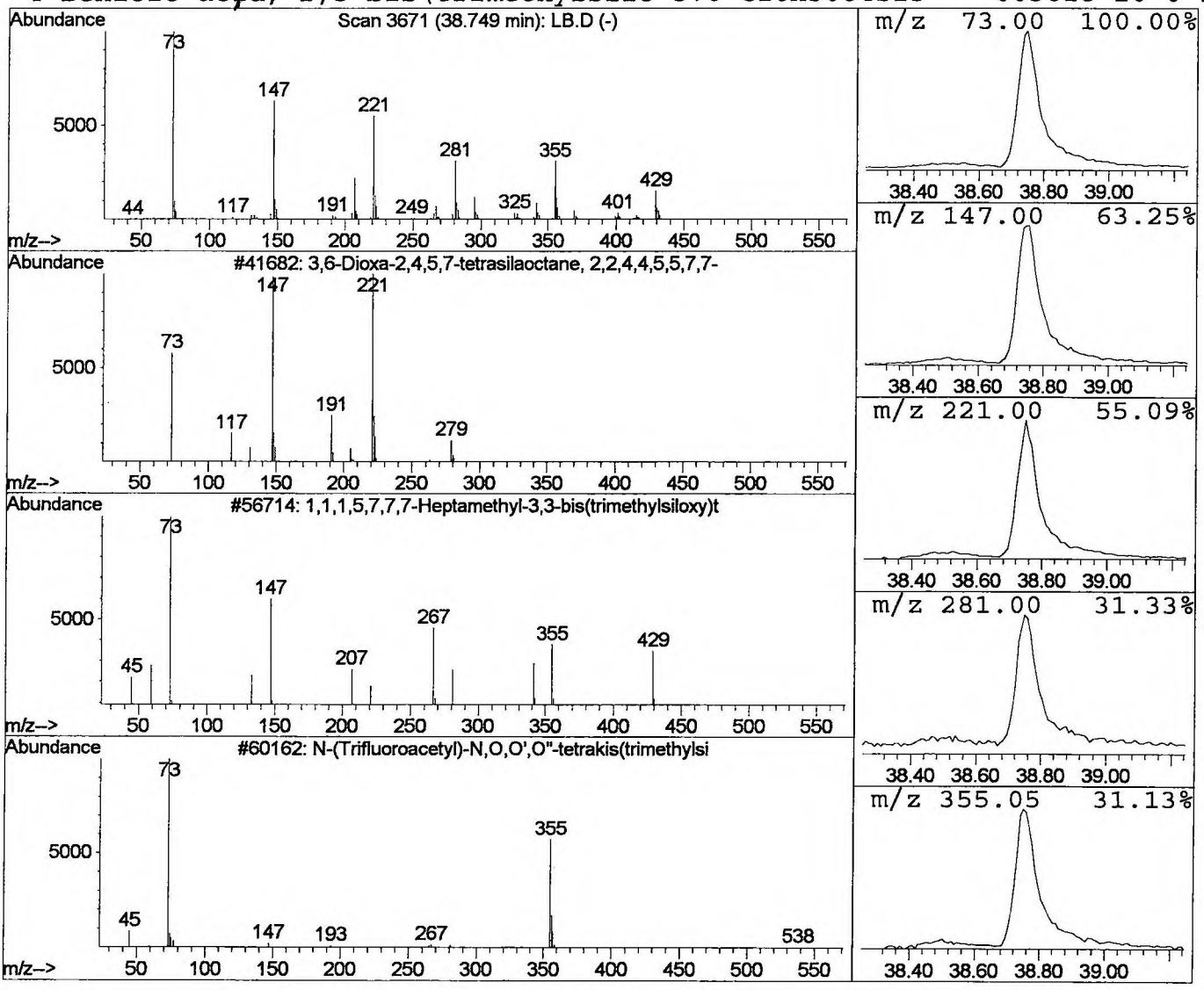
Quant Method : C:\HPCHEM\1\METHODS\GCMS0412.M (RTE Integrator)
Title : 209 625/TCLP calibration table
Library : C:\DATABASE\NBS75K.L

****

Peak Number 6 3,6-Dioxa-2,4,5,7-tetrasilaoct Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
38.75	8.80 ug/l	788207	Perylene-d12	37.87

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	3,6-Dioxa-2,4,5,7-tetrasilaoctane,	294	C10H30O2Si4	004342-25-0	32
2	1,1,1,5,7,7,7-Heptamethyl-3,3-bis(t	444	C13H40O5Si6	038147-00-1	28
3	N-(Trifluoroacetyl)-N,O,O',O''-tetr	553	C22H42F3NO4Si4	000000-00-0	10
4	Benzoic acid, 2,5-bis(trimethylsilo	370	C16H30O4Si3	003618-20-0	10



Library Search Compound Report

Data File : C:\HPCHEM\1\DATA\APR1702\LB.D Vial: 3
Acq On : 17 Apr 2002 10:13 am Operator:
Sample : gc/ms scan 3/19 Inst : 209
Misc : Multiplr: 1.00
MS Integration Params: LSCINT.P

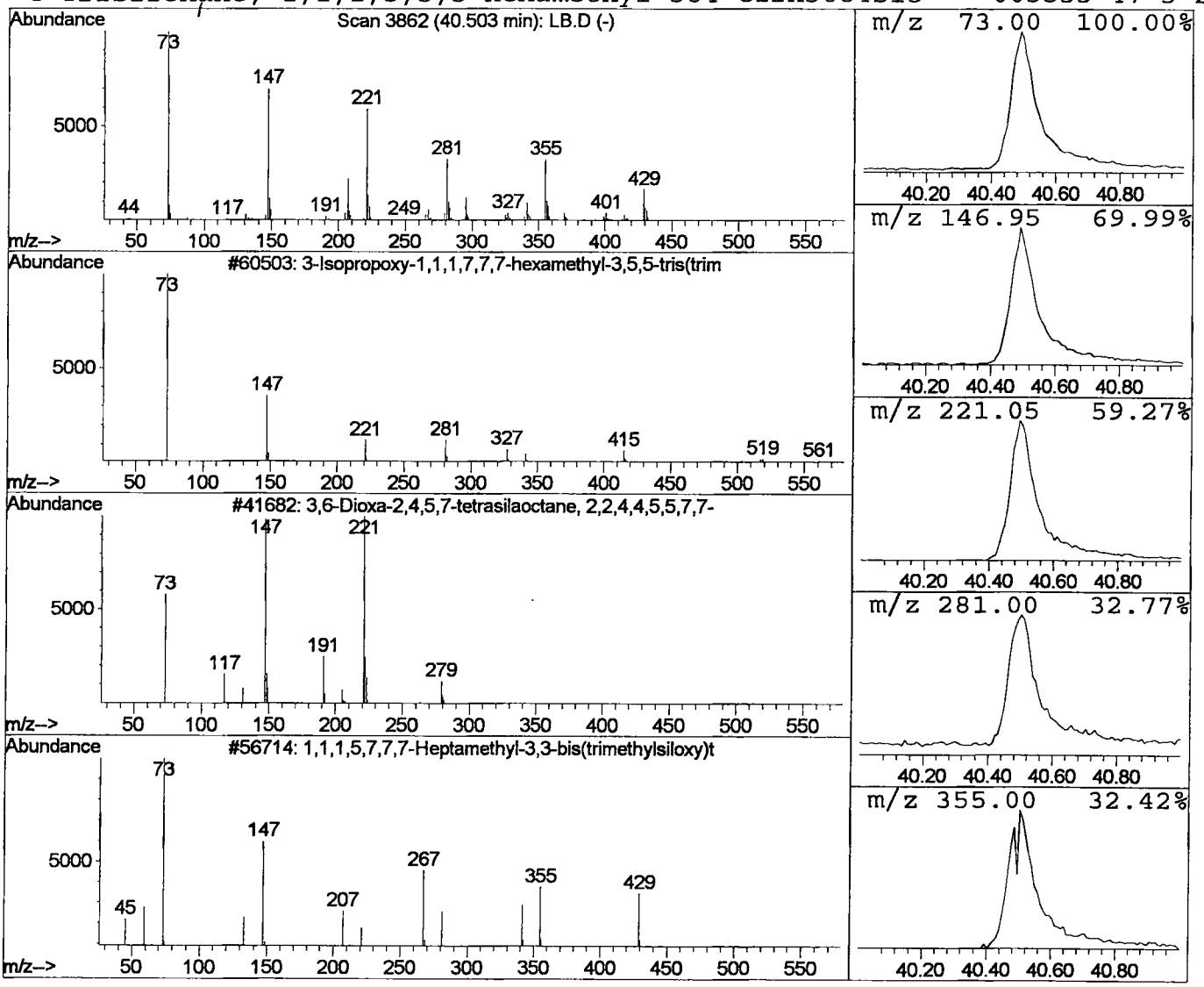
Quant Method : C:\HPCHEM\1\METHODS\GCMS0412.M (RTE Integrator)
Title : 209 625/TCLP calibration table
Library : C:\DATABASE\NBS75K.L

****

Peak Number 7 3-Isopropoxy-1,1,1,7,7,7-hexam Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
40.50	8.80 ug/l	788043	Perylene-d12	37.87

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	3-Isopropoxy-1,1,1,7,7,7-hexamethyl	576	C18H52O7Si7	071579-69-6	37
2	3,6-Dioxa-2,4,5,7-tetrasilaoctane,	294	C10H30O2Si4	004342-25-0	32
3	1,1,1,5,7,7,7-Heptamethyl-3,3-bis(t	444	C13H40O5Si6	038147-00-1	28
4	Trisiloxane, 1,1,1,5,5,5-hexamethyl	384	C12H36O4Si5	003555-47-3	25



Library Search Compound Report

Data File : C:\HPCHEM\1\DATA\APR1702\LB.D Vial: 3
Acq On : 17 Apr 2002 10:13 am Operator:
Sample : gc/ms scan 3/19 Inst : 209
Misc : Multiplr: 1.00
MS Integration Params: LSCINT.P

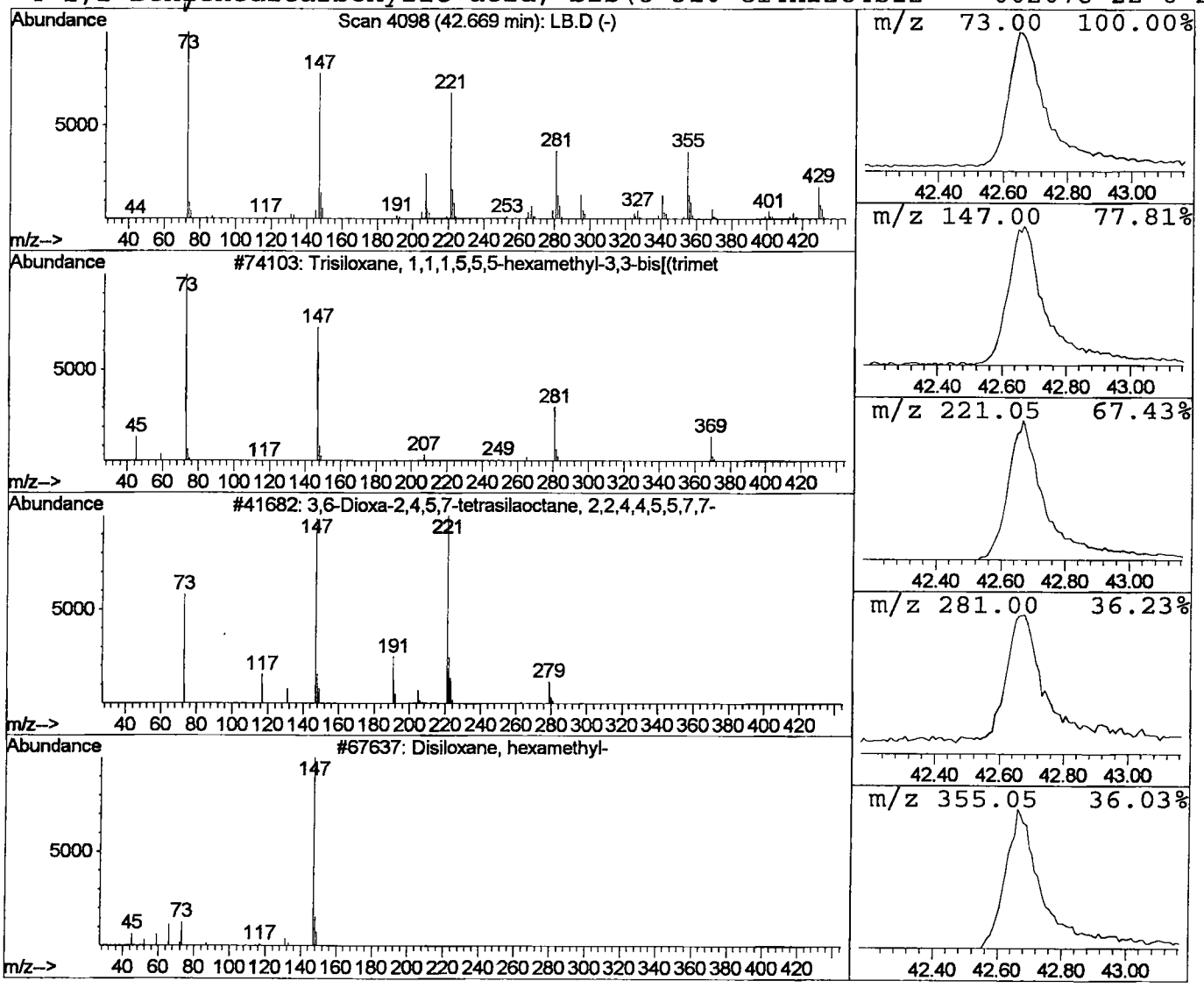
Quant Method : C:\HPCHEM\1\METHODS\GCMS0412.M (RTE Integrator)
Title : 209 625/TCLP calibration table
Library : C:\DATABASE\NBS75K.L

****

Peak Number 8 Trisiloxane, 1,1,1,5,5,5-hexam Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
42.67	7.28 ug/l	651983	Perylene-d12	37.87

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Trisiloxane, 1,1,1,5,5,5-hexamethyl	384	C12H36O4Si5	003555-47-3	43
2	3,6-Dioxa-2,4,5,7-tetrasilaoctane,	294	C10H30O2Si4	004342-25-0	28
3	Disiloxane, hexamethyl-	162	C6H18OSi2	000107-46-0	27
4	1,2-Benzenedicarboxylic acid, bis(t	310	C14H22O4Si2	002078-22-0	27



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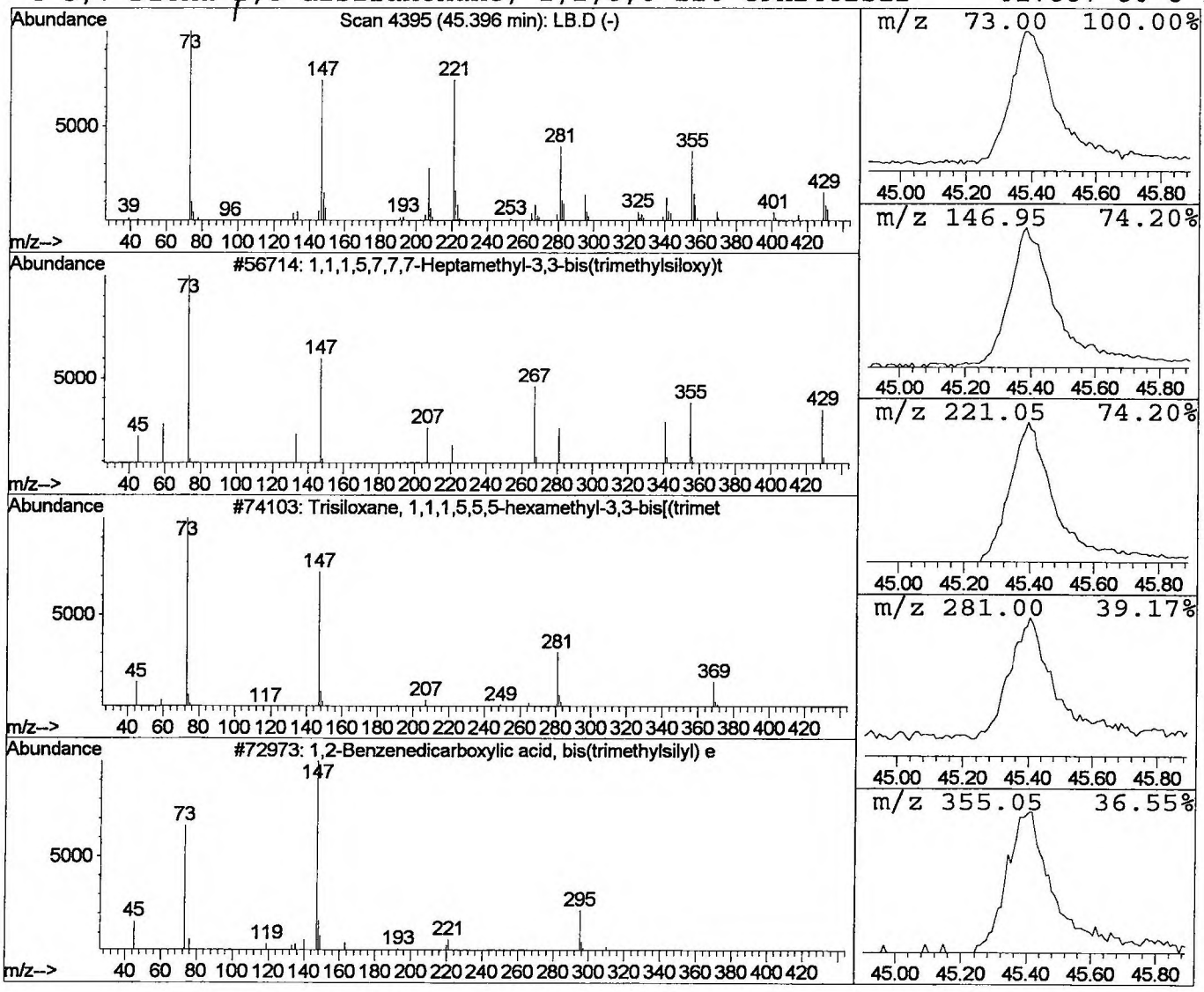
****

Peak Number 9 1,1,1,5,7,7,7-Heptamethyl-3,3- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
45.40	5.57 ug/l	498841	Perylene-d12	37.87

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	Trisiloxane, 1,1,1,5,5,5-hexamethyl	384	C12H36O4Si5	003555-47-3	25
3	1,2-Benzenedicarboxylic acid, bis(t	310	C14H22O4Si2	002078-22-0	14
4	3,7-Dioxa-2,8-disilanonane, 2,2,8,8	220	C9H24O2Si2	017887-80-8	14



Tentatively Identified Compound (LSC) summary

Operator ID: Date Acquired: 17 Apr 2002 10:13 am
Data File: C:\HPCHEM\1\DATA\APR1702\LB.D
Name: gc/ms scan 3/19
Misc:
Method: C:\HPCHEM\1\METHODS\GCMS0412.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
N-Ethylformamide	8.11	4.3	ug/l	578215	ISTD01	12.21	2674290	20.0
Cyclopentasiloxane,	14.82	0.2	ug/l	40517	ISTD02	15.84	3372190	20.0
3-Trimethylsilyloxys	34.89	1.0	ug/l	111628	ISTD05	33.64	2335020	20.0
1,2-Benzenedicarboxy	36.09	3.1	ug/l	278661	ISTD06	37.87	1791450	20.0
3-Isopropoxy-1,1,1,7	37.31	6.0	ug/l	537375	ISTD06	37.87	1791450	20.0
3,6-Dioxa-2,4,5,7-te	38.75	8.8	ug/l	788207	ISTD06	37.87	1791450	20.0
3-Isopropoxy-1,1,1,7	40.50	8.8	ug/l	788043	ISTD06	37.87	1791450	20.0
Trisiloxane, 1,1,1,5	42.67	7.3	ug/l	651983	ISTD06	37.87	1791450	20.0
1,1,1,5,5,7,7-Heptam	45.40	5.6	ug/l	498841	ISTD06	37.87	1791450	20.0

column

LB.D GCMS0412.M Thu Apr 18 10:56:38 2002