& Research
Time: 16:20:29

Sample: AE02581
Analysis: $B1GCMS Blank for GCMS Scan
Units: ug
Started: 3/13/02 16:16 Ended: 3/13/02 16:16 Analyst: MG
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	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/1,2-Diphenylhydraz		< 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\MAR0702\LB.D Vial: 15
Acq On : 8 Mar 2002 3:29 am Operator:
Sample : gc/ms scan 2/5 Inst : GC/MS 209
Misc : Multiplr: 1.00
MS Integration Params: GOOFY.P
Quant Time: Mar 13 14:21 2002 Quant Results File: GCMS0208.RES

Quant Method : C:\HPCHEM\1\METHODS\GCMS0208.M (RTE Integrator)
Title : 209 625/TCLP calibration table
Last Update : Wed Feb 27 11:30:15 2002
Response via : Initial Calibration
DataAcq Meth : GCMS0208

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	12.30	152	410885	20.00	ug/l	-0.05
10) Naphthalene-d8	15.94	136	1579826	20.00	ug/l	-0.06
17) Acenaphthene-d10	21.25	164	844719	20.00	ug/l	-0.06
29) Phenanthrene-d10	25.69	188	1293996	20.00	ug/l	-0.07
38) Chrysene-d12	33.76	240	1051389	20.00	ug/l	-0.08
44) Perylene-d12	38.02	264	783741	20.00	ug/l	-0.10
System Monitoring Compounds						
37) 2,4-DDD	30.77	235	72731	5.38	ug/mL	0.27
Spiked Amount 5.000			Recovery = 107.60%			
Target Compounds						Qvalue
2) N-nitroso-dimethyl amine	0.00	74	0	N.D.		
3) bis(2-Chloroethyl)ether	0.00	93	0	N.D.		
4) 1,3-Dichlorobenzene	0.00	146	0	N.D.		
5) 1,4-Dichlorobenzene	0.00	146	0	N.D.		
6) 1,2-Dichlorobenzene	0.00	146	0	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	0.00	77	0	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	15.99	128	524	N.D.		
16) Hexachlorobutadiene	0.00	225	0	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	21.65	165	229	N.D.		
25) Diethylphthalate	0.00	149	0	N.D.		
26) 4-Chlorophenyl-phenylether	0.00	204	0	N.D.		
27) Fluorene	0.00	166	0	N.D.		
28) Azobenzen/ 1,2-Diphenylhyd	0.00	77	0	N.D.		

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

Quantitation Report (QT Reviewed)

Data File : C:\HPCHEM\1\DATA\MAR0702\LB.D Vial: 15
Acq On : 8 Mar 2002 3:29 am Operator:
Sample : gc/ms scan 2/5 Inst : GC/MS 209
Misc : Multiplr: 1.00
MS Integration Params: GOOFY.P
Quant Time: Mar 13 14:21 2002 Quant Results File: GCMS0208.RES

Quant Method : C:\HPCHEM\1\METHODS\GCMS0208.M (RTE Integrator)
Title : 209 625/TCLP calibration table
Last Update : Wed Feb 27 11:30:15 2002
Response via : Initial Calibration
DataAcq Meth : GCMS0208

	Compound	R.T.	QIon	Response	Conc Unit	Qvalue
30)	N-Nitrosodiphenylamine	0.00	168	0	N.D.	
31)	4-Bromophenyl-phenylether	0.00	248	0	N.D.	
32)	Hexachlorobenzene	0.00	284	0	N.D.	
33)	Phenanthrene	25.69	178	364	N.D.	
34)	Anthracene	25.69	178	364	N.D.	
35)	Di-n-butylphthalate	27.66	149	20245	N.D.	
36)	Fluoranthene	0.00	202	0	N.D.	
39)	Pyrene	0.00	202	0	N.D.	
40)	Butylbenzylphthalate	32.35	149	187	N.D.	
41)	Benzo(a)anthracene	33.76	228	2492	N.D.	
42)	Bis(2-ethylhexyl)phthalate	33.97	149	2909	N.D.	
43)	Chrysene	33.76	228	2491	N.D.	
45)	Di-n-Octylphthalate	36.17	149	1038	N.D.	
46)	Benzo(b)fluoranthene	0.00	252	0	N.D.	
47)	Benzo(k)fluoranthene	0.00	252	0	N.D.	
48)	Benzo(a)pyrene	38.02	252	3172	N.D.	
49)	Indeno(1,2,3-cd)pyrene	0.00	276	0	N.D.	
50)	Dibenz(a,h)anthracene	0.00	278	0	N.D.	
51)	Benzo(g,h,i)perylene	0.00	276	0	N.D.	

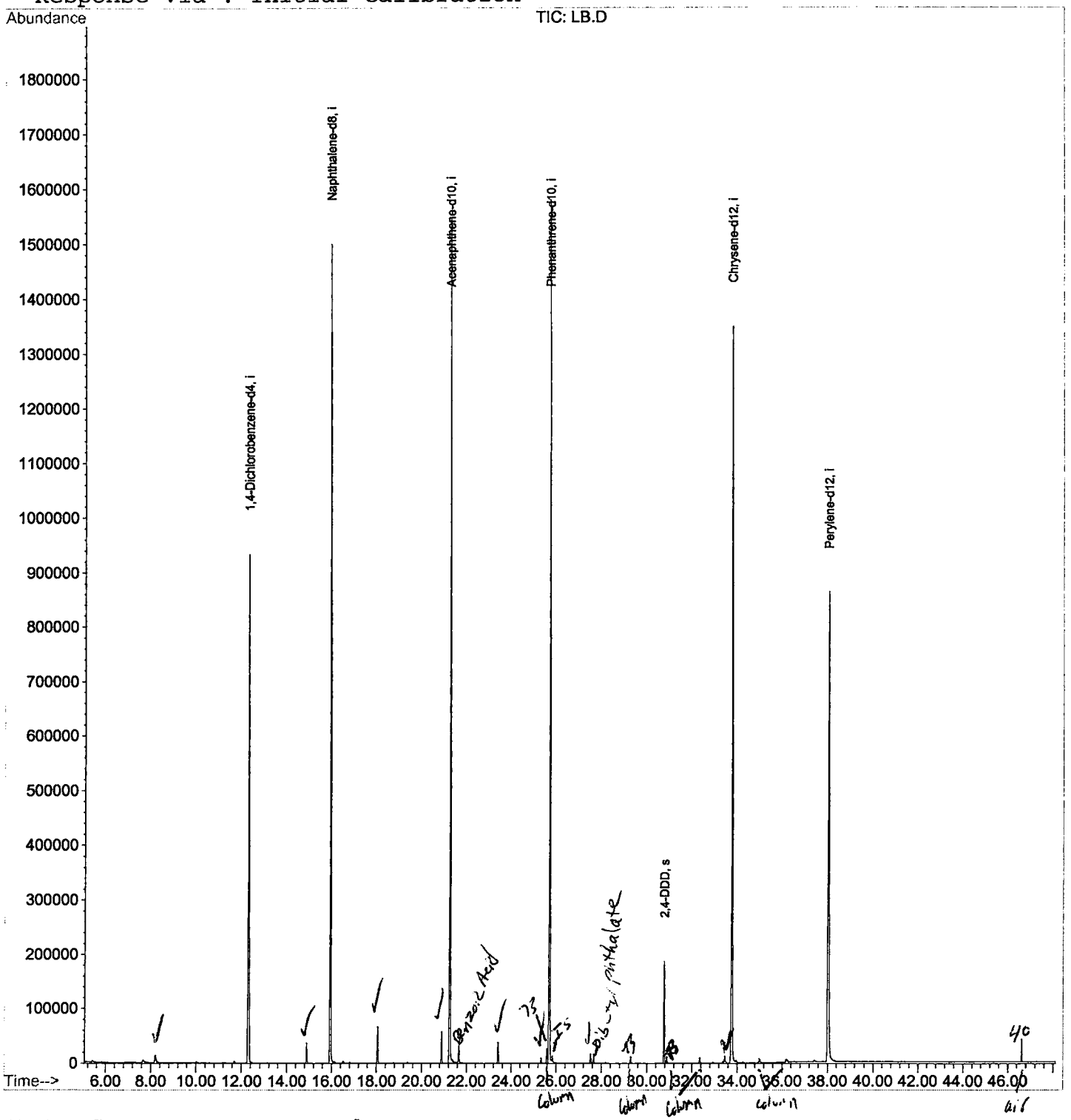
(#) = qualifier out of range (m) = manual integration

Quantitation Report

Data File : C:\HPCHEM\1\DATA\MAR0702\LB.D
Acq On : 8 Mar 2002 3:29 am
Sample : gc/ms scan 2/5
Misc :
MS Integration Params: GOOFY.P
Quant Time: Mar 13 14:21 2002

Vial: 15
Operator:
Inst : GC/MS 209
Multiplr: 1.00
Quant Results File: GCMS0208.RES

Method : C:\HPCHEM\1\METHODS\GCMS0208.M (RTE Integrator)
Title : 209 625/TCLP calibration table
Last Update : Wed Feb 27 11:30:15 2002
Response via : Initial Calibration



LSC Area Percent Report

Data File : C:\HPCHEM\1\DATA\MAR0702\LB.D Vial: 15
Acq On : 8 Mar 2002 3:29 am Operator:
Sample : gc/ms scan 2/5 Inst : GC/MS 209
Misc : Multiplr: 1.00
MS Integration Params: LSCINT.P

Method : C:\HPCHEM\1\METHODS\GCMS0208.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

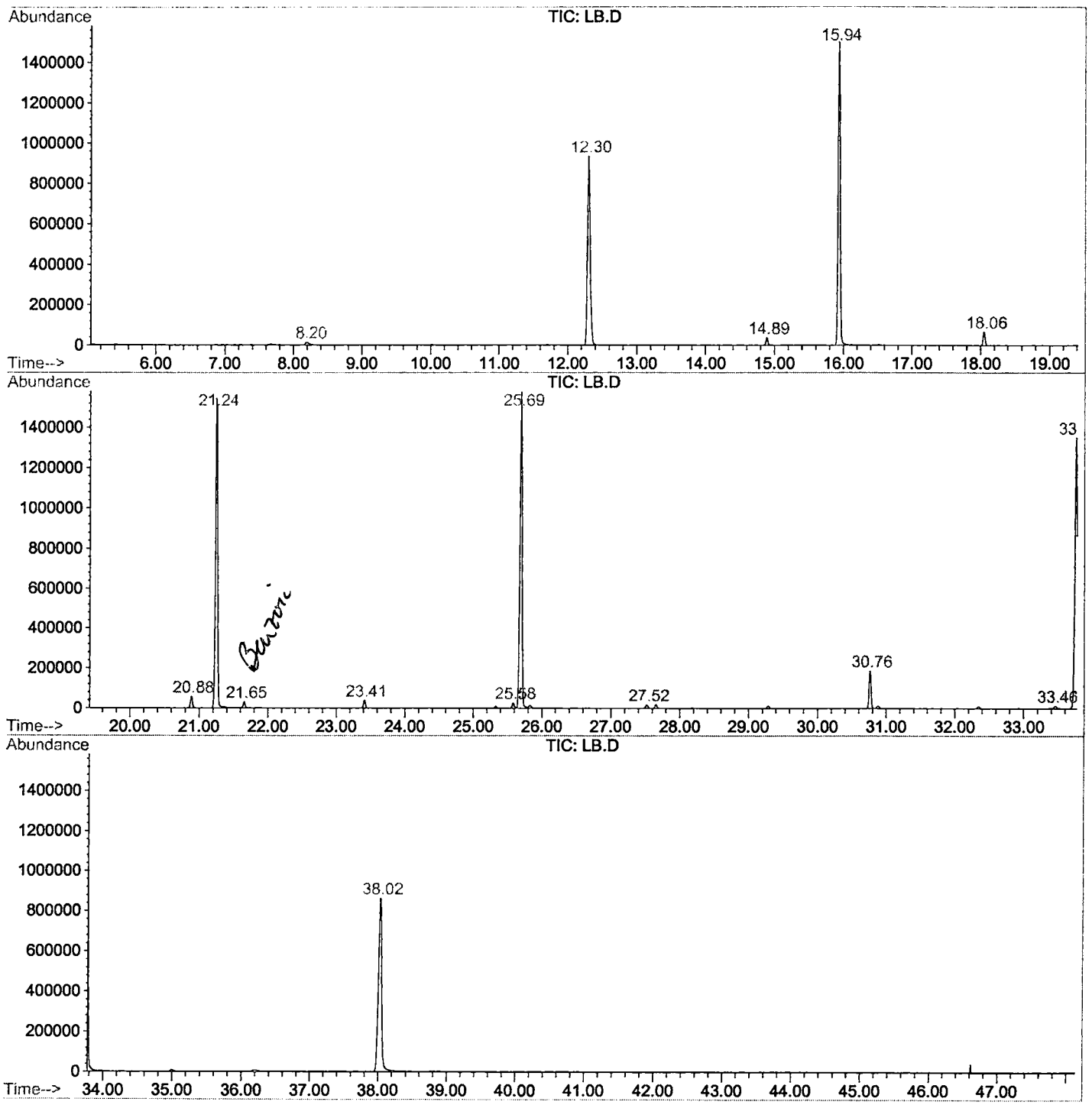
peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	peak area	peak % max.	% of total
1	8.197	337	343	356	rBV3	13831	53325	1.58%	0.281%
2	12.300	784	790	804	rBV	932743	2300974	68.08%	12.108%
3	14.889	1067	1072	1077	rBV	37137	71158	2.11%	0.374%
4	15.936	1180	1186	1201	rBV	1499726	3046503	90.14%	16.031%
5	18.056	1411	1417	1425	rVB	66558	132290	3.91%	0.696%
6	20.884	1719	1725	1733	rVB	57547	116503	3.45%	0.613%
7	21.242	1758	1764	1774	rBV	1544525	3293413	97.45%	17.331%
8	21.655	1804	1809	1814	rVB	30102	52141	1.54%	0.274%
9	23.408	1994	2000	2005	rBV	38132	77217	2.28%	0.406%
10	25.584	2232	2237	2242	rBV	26823	51181	1.51%	0.269%
11	25.694	2242	2249	2259	rBV2	1579859	3379723	100.00%	17.785%
12	27.521	2443	2448	2453	rVB	17274	33925	1.00%	0.179%
13	30.762	2796	2801	2808	rBV	187196	375203	11.10%	1.974%
14	33.461	3089	3095	3106	rVB	12365	34987	1.04%	0.184%
15	33.764	3116	3128	3142	rBV2	1349347	3202653	94.76%	16.853%
16	38.023	3583	3592	3610	rBV2	860985	2782206	82.32%	14.641%

Sum of corrected areas: 19003402

LB.D GCMS0208.M Wed Mar 13 15:03:21 2002

LSC Report - Integrated Chromatogram

File : C:\HPCHEM\1\DATA\MAR0702\LB.D
Operator :
Acquired : 8 Mar 2002 3:29 am using AcqMethod GCMS0208
Instrument : GC/MS 209
Sample Name: gc/ms scan 2/5
Misc Info :
Vial Number: 15
Quant File :GCMS0208.RES (RTE Integrator)



LB.D GCMS0208.M Wed Mar 13 15:03:26 2002

Library Search Compound Report

Data File : C:\HPCHEM\1\DATA\MAR0702\LB.D Vial: 15
Acq On : 8 Mar 2002 3:29 am Operator:
Sample : gc/ms scan 2/5 Inst : GC/MS 209
Misc : Multiplr: 1.00
MS Integration Params: LSCINT.P

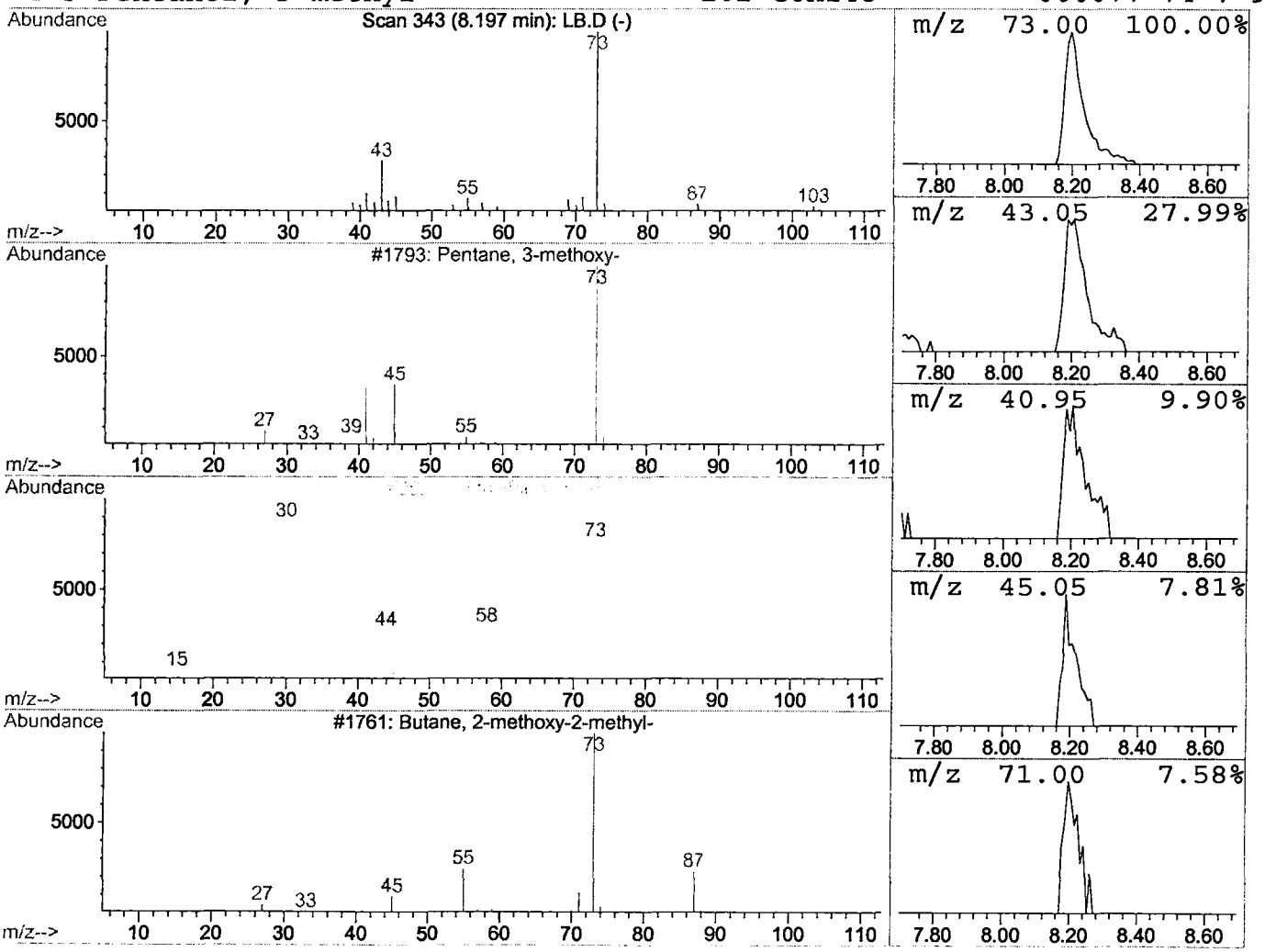
Quant Method : C:\HPCHEM\1\METHODS\GCMS0208.M (RTE Integrator)
Title : 209 625/TCLP calibration table
Library : C:\DATABASE\NBS75K.L

****

Peak Number 1 Pentane, 3-methoxy- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.20	0.46 ug/l	53325	1,4-Dichlorobenzene-d4	12.30

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Pentane, 3-methoxy-	102	C6H14O	036839-67-5	28
2	N-Ethylformamide	73	C3H7NO	000627-45-2	9
3	Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	9
4	3-Pentanol, 3-methyl-	102	C6H14O	000077-74-7	9



Library Search Compound Report

Data File : C:\HPCHEM\1\DATA\MAR0702\LB.D Vial: 15
Acq On : 8 Mar 2002 3:29 am Operator:
Sample : gc/ms scan 2/5 Inst : GC/MS 209
Misc : Multiplr: 1.00
MS Integration Params: LSCINT.P

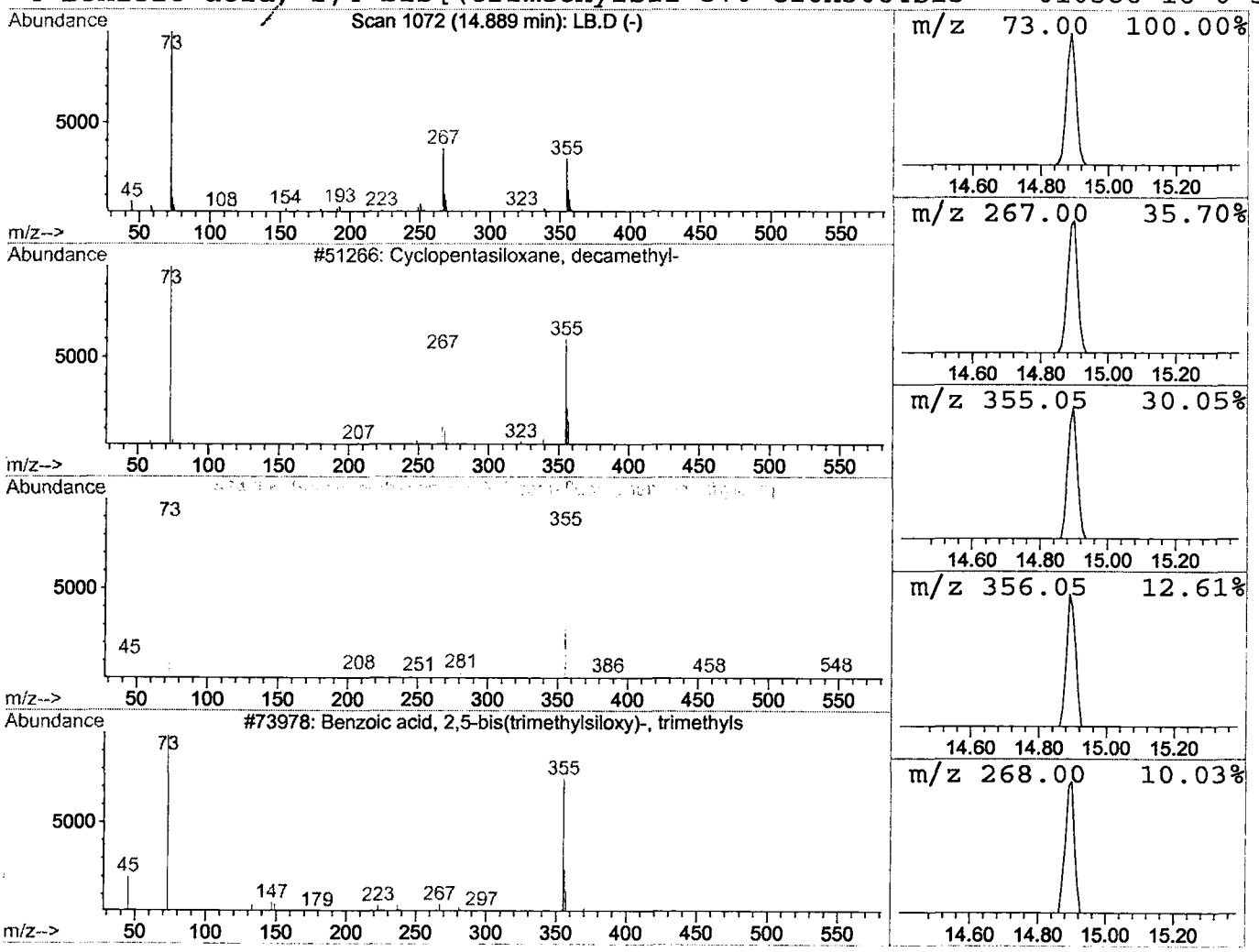
Quant Method : C:\HPCHEM\1\METHODS\GCMS0208.M (RTE Integrator)
Title : 209 625/TCLP calibration table
Library : C:\DATABASE\NBS75K.L

****

Peak Number 2 Cyclopentasiloxane, decamethyl Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.89	0.47 ug/l	71158	Naphthalene-d8	15.94

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclopentasiloxane, decamethyl-	370	C10H30O5Si5	000541-02-6	83
2	Benzeneethanamine, N-[(pentafluorop	563	C24H34F5NO3Si3	055429-13-5	43
3	Benzoic acid, 2,5-bis(trimethylsilo	370	C16H30O4Si3	003618-20-0	37
4	Benzoic acid, 2,4-bis[(trimethylsil	370	C16H30O4Si3	010586-16-0	37



Library Search Compound Report

Data File : C:\HPCHEM\1\DATA\MAR0702\LB.D Vial: 15
Acq On : 8 Mar 2002 3:29 am Operator:
Sample : gc/ms scan 2/5 Inst : GC/MS 209
Misc : Multiplr: 1.00
MS Integration Params: LSCINT.P

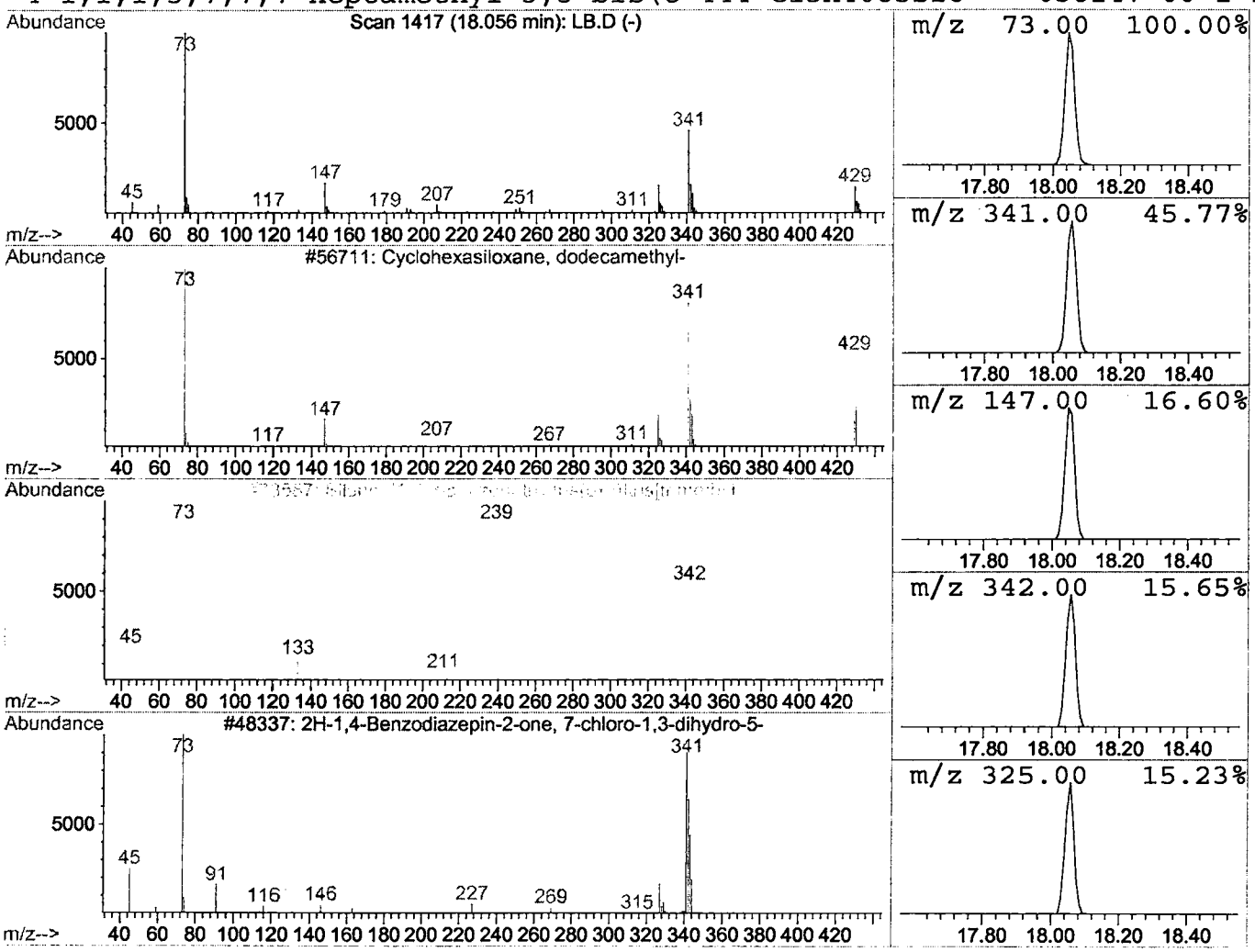
Quant Method : C:\HPCHEM\1\METHODS\GCMS0208.M (RTE Integrator)
Title : 209 625/TCLP calibration table
Library : C:\DATABASE\NBS75K.L

****

Peak Number 3 Cyclohexasiloxane, dodecamethy Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.06	0.87 ug/l	132290	Naphthalene-d8	15.94

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclohexasiloxane, dodecamethyl-	444	C12H36O6Si6	000540-97-6	59
2	Silane, [1,2,3-benzenetriyltris(oxy	342	C15H30O3Si3	017864-23-2	22
3	2H-1,4-Benzodiazepin-2-one, 7-chlor	342	C18H19ClN2OSi	055299-24-6	22
4	1,1,1,5,7,7,7-Heptamethyl-3,3-bis(t	444	C13H40O5Si6	038147-00-1	12



Library Search Compound Report

Data File : C:\HPCHEM\1\DATA\MAR0702\LB.D
Acq On : 8 Mar 2002 3:29 am
Sample : gc/ms scan 2/5
Misc :
MS Integration Params: LSCINT.P

Vial: 15
Operator:
Inst : GC/MS 209
Multiplr: 1.00

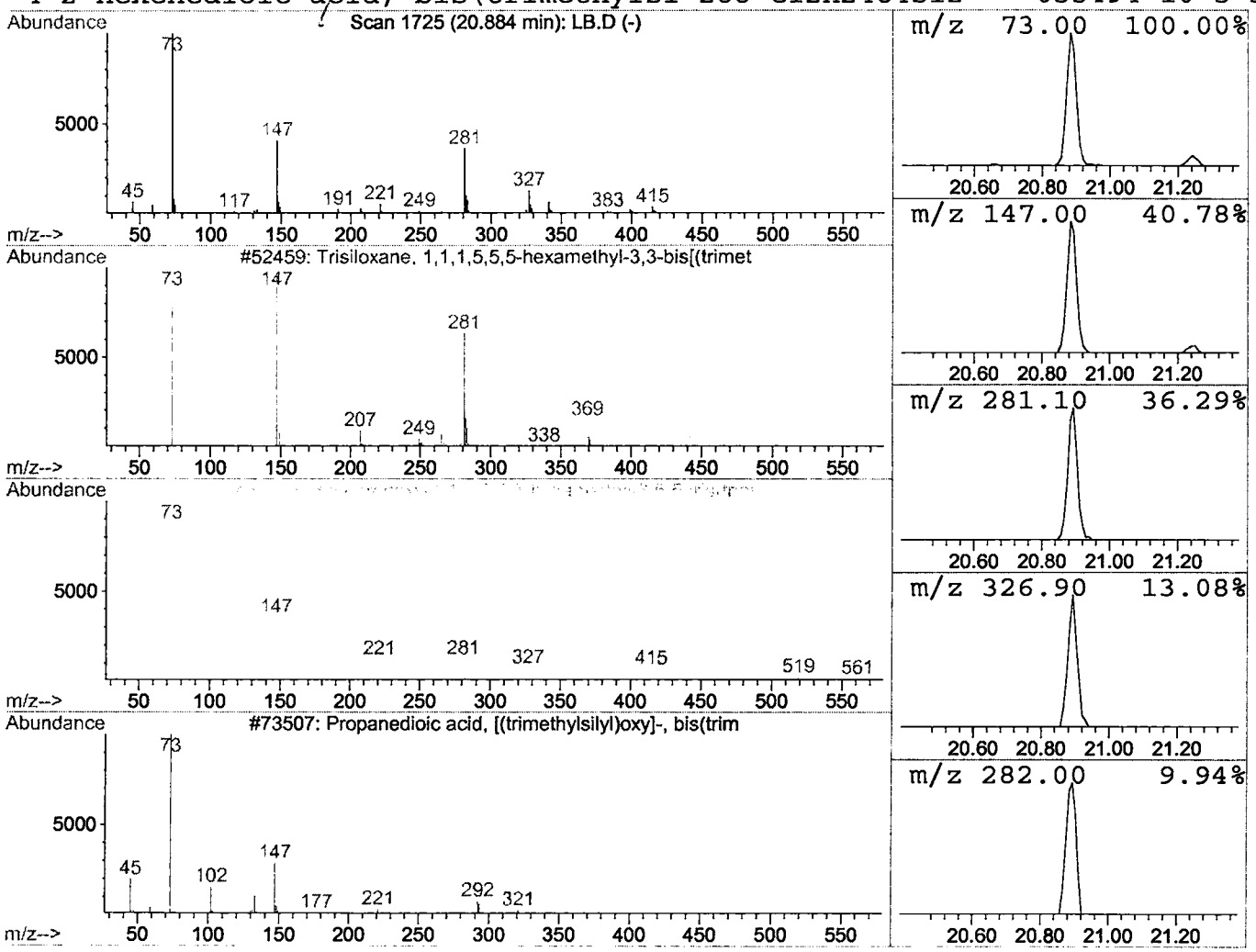
Quant Method : C:\HPCHEM\1\METHODS\GCMS0208.M (RTE Integrator)
Title : 209 625/TCLP calibration table
Library : C:\DATABASE\NBS75K.L

****

Peak Number 4 Trisiloxane, 1,1,1,5,5,5-hexam Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.88	0.71 ug/l	116503	Acenaphthene-d10	21.25

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Trisiloxane, 1,1,1,5,5,5-hexamethyl	384	C12H36O4Si5	003555-47-3	45
2	3-Isopropoxy-1,1,1,7,7,7-hexamethyl	576	C18H52O7Si7	071579-69-6	45
3	Propanedioic acid, [(trimethylsilyl	336	C12H28O5Si3	038165-93-4	38
4	2-Hexenedioic acid, bis(trimethylsi	288	C12H24O4Si2	055494-10-5	32



Library Search Compound Report

Data File : C:\HPCHEM\1\DATA\MAR0702\LB.D Vial: 15
Acq On : 8 Mar 2002 3:29 am Operator:
Sample : gc/ms scan 2/5 Inst : GC/MS 209
Misc : Multiplr: 1.00
MS Integration Params: LSCINT.P

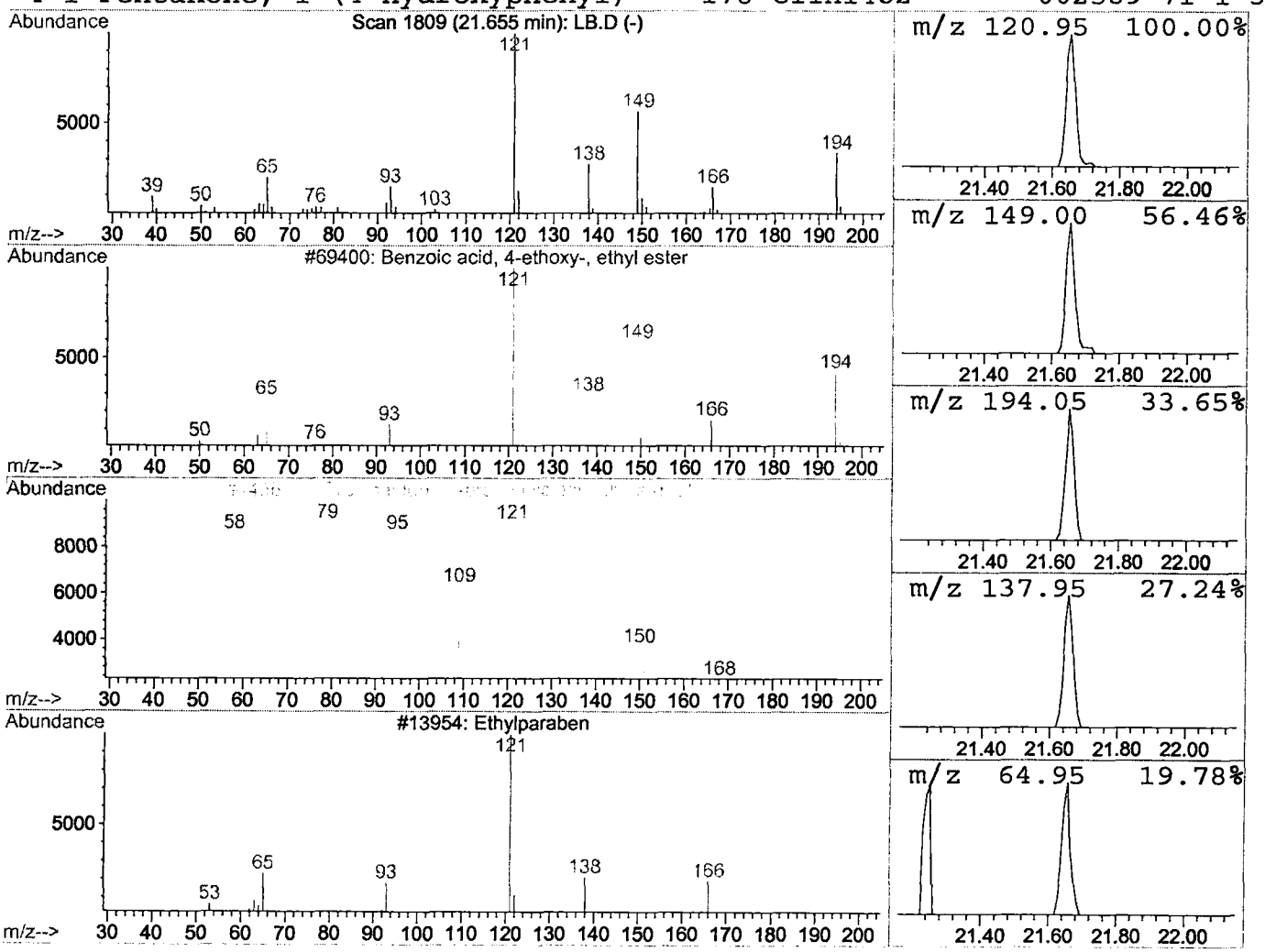
Quant Method : C:\HPCHEM\1\METHODS\GCMS0208.M (RTE Integrator)
Title : 209 625/TCLP calibration table
Library : C:\DATABASE\NBS75K.L

****

Peak Number 5 Benzoic acid, 4-ethoxy-, ethyl Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.65	0.32 ug/l	52141	Acenaphthene-d10	21.25

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzoic acid, 4-ethoxy-, ethyl este	194	C11H14O3	023676-09-7	97
2	1-Cyclopentene-1-propanoic acid, .b	168	C10H16O2	061833-30-5	80
3	Ethylparaben	166	C9H10O3	000120-47-8	60
4	1-Pentanone, 1-(4-hydroxyphenyl)-	178	C11H14O2	002589-71-1	38



Library Search Compound Report

Data File : C:\HPCHEM\1\DATA\MAR0702\LB.D
Acq On : 8 Mar 2002 3:29 am
Sample : gc/ms scan 2/5
Misc :
MS Integration Params: LSCINT.P

Vial: 15
Operator:
Inst : GC/MS 209
Multiplr: 1.00

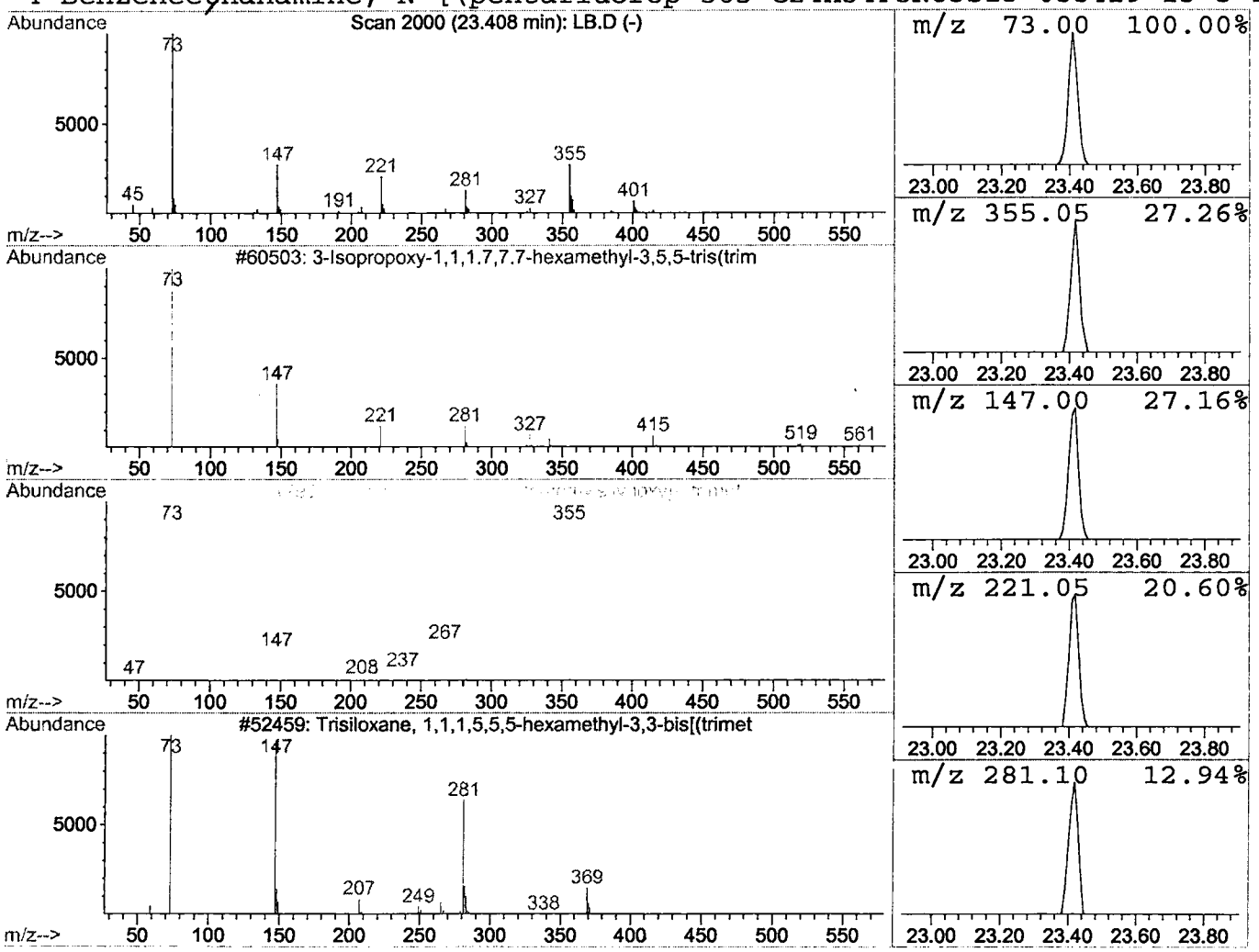
Quant Method : C:\HPCHEM\1\METHODS\GCMS0208.M (RTE Integrator)
Title : 209 625/TCLP calibration table
Library : C:\DATABASE\NBS75K.L

****

Peak Number 6 3-Isopropoxy-1,1,1,7,7,7-hexam Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.41	0.47 ug/l	77217	Acenaphthene-d10	21.25

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	Benzoic acid, 2,6-bis[(trimethylsil	370	C16H30O4Si3	003782-85-2	23
3	Trisiloxane, 1,1,1,5,5,5-hexamethyl	384	C12H36O4Si5	003555-47-3	17
4	Benzeneethanamine, N-[(pentafluorop	563	C24H34F5NO3Si3	055429-13-5	17



Library Search Compound Report

Data File : C:\HPCHEM\1\DATA\MAR0702\LB.D Vial: 15
Acq On : 8 Mar 2002 3:29 am Operator:
Sample : gc/ms scan 2/5 Inst : GC/MS 209
Misc : Multiplr: 1.00
MS Integration Params: LSCINT.P

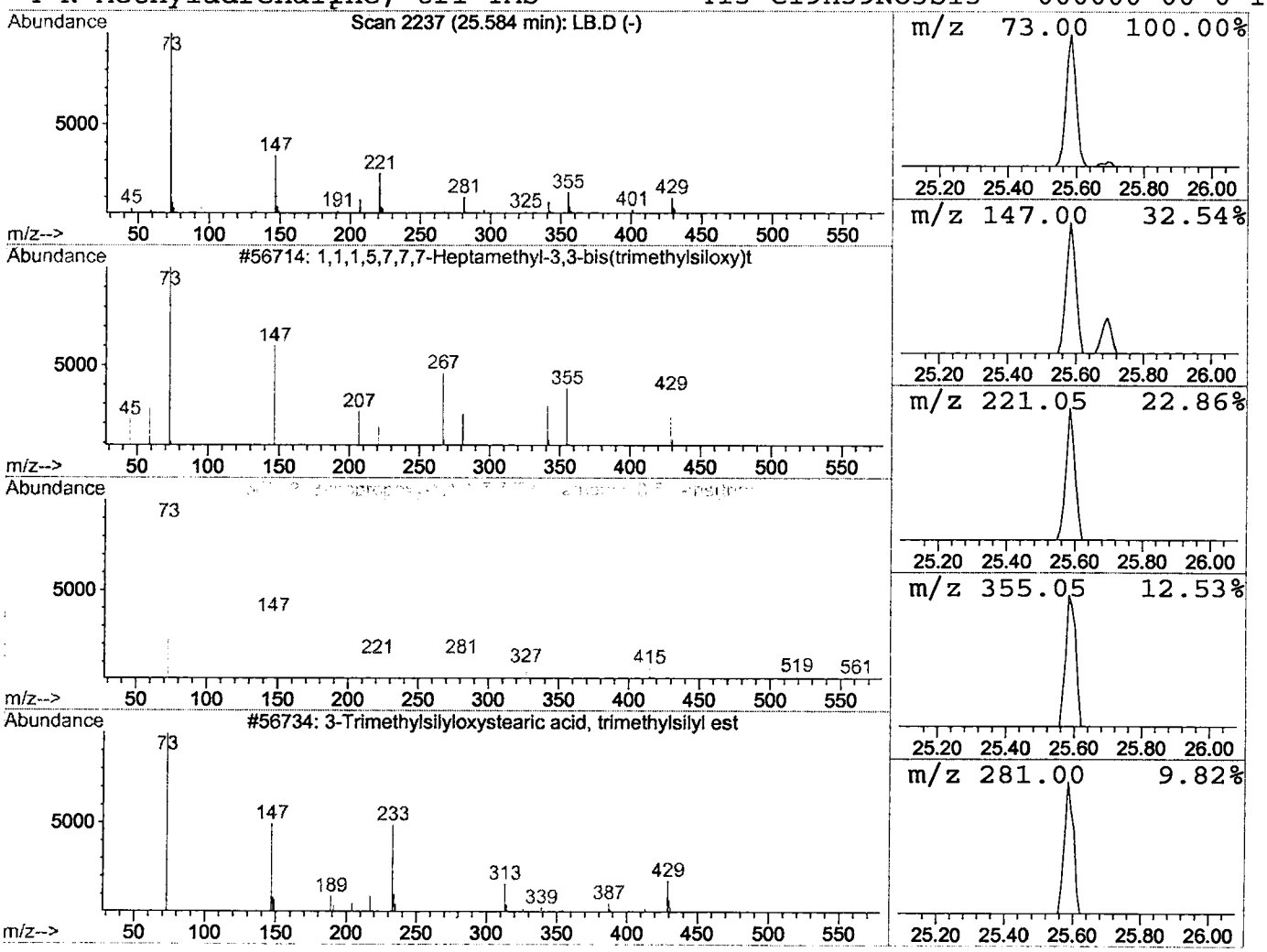
Quant Method : C:\HPCHEM\1\METHODS\GCMS0208.M (RTE Integrator)
Title : 209 625/TCLP calibration table
Library : C:\DATABASE\NBS75K.L

****

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
25.58	0.30 ug/l	51181	Phenanthrene-d10	25.69

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	47
2	3-Isopropoxy-1,1,1,7,7,7-hexamethyl	576	C18H52O7Si7	071579-69-6	25
3	3-Trimethylsilyloxystearic acid, tr	444	C24H52O3Si2	000000-00-0	17
4	N-Methyladrenaline, tri-TMS	413	C19H39NO3Si3	000000-00-0	12



Library Search Compound Report

Data File : C:\HPCHEM\1\DATA\MAR0702\LB.D Vial: 15
Acq On : 8 Mar 2002 3:29 am Operator:
Sample : gc/ms scan 2/5 Inst : GC/MS 209
Misc : Multiplr: 1.00
MS Integration Params: LSCINT.P

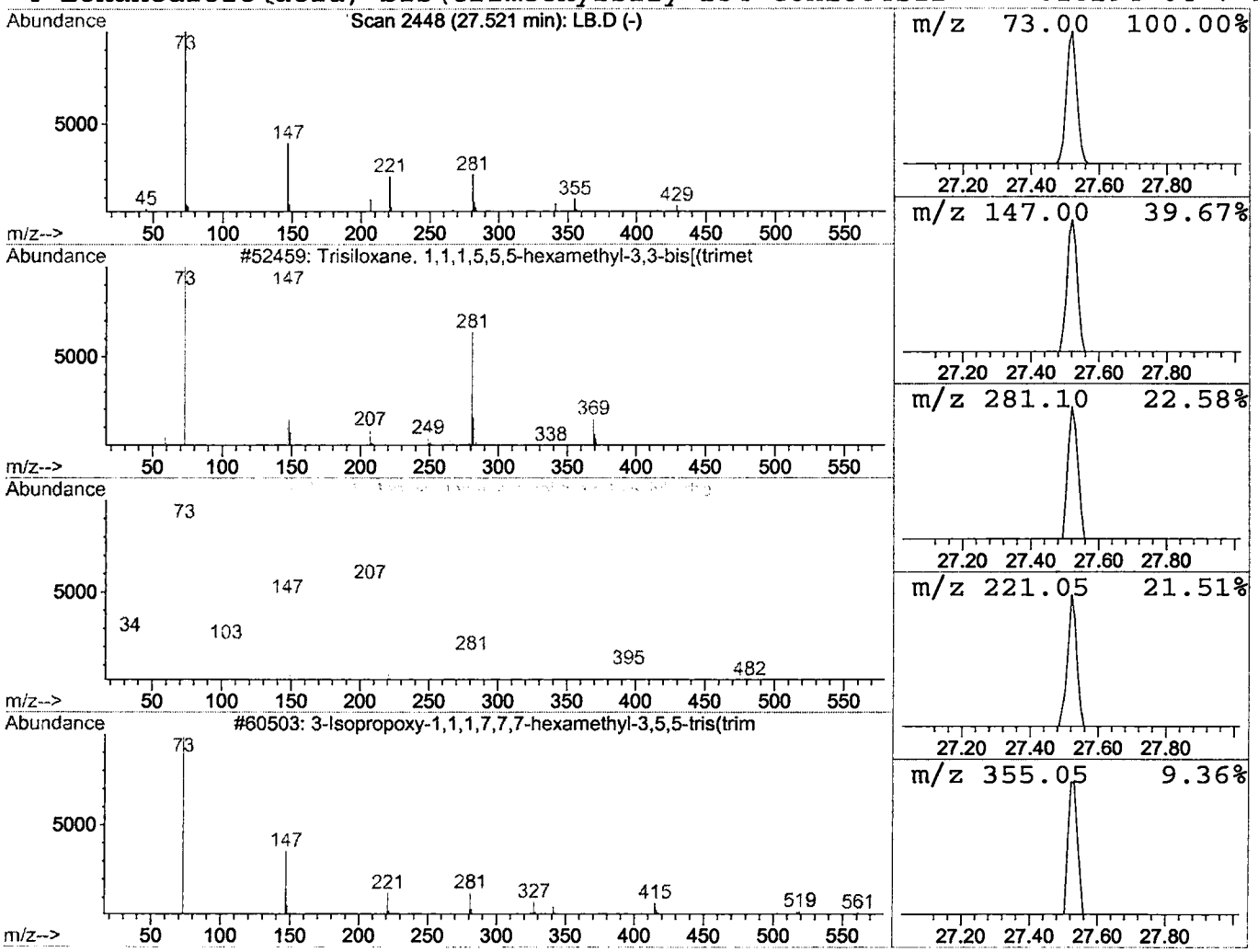
Quant Method : C:\HPCHEM\1\METHODS\GCMS0208.M (RTE Integrator)
Title : 209 625/TCLP calibration table
Library : C:\DATABASE\NBS75K.L

****

Peak Number 8 Trisiloxane, 1,1,1,5,5,5-hexam Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
27.52	0.20 ug/l	33925	Phenanthrene-d10	25.69

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Trisiloxane, 1,1,1,5,5,5-hexamethyl	384	C12H36O4Si5	003555-47-3	50
2	1-Monolinoleoylglycerol trimethylsi	498	C27H54O4Si2	054284-45-6	36
3	3-Isopropoxy-1,1,1,7,7,7-hexamethyl	576	C18H52O7Si7	071579-69-6	33
4	Ethanedioic acid, bis(trimethylsily	234	C8H18O4Si2	018294-04-7	17



Library Search Compound Report

Data File : C:\HPCHEM\1\DATA\MAR0702\LB.D Vial: 15
Acq On : 8 Mar 2002 3:29 am Operator:
Sample : gc/ms scan 2/5 Inst : GC/MS 209
Misc : Multiplr: 1.00
MS Integration Params: LSCINT.P

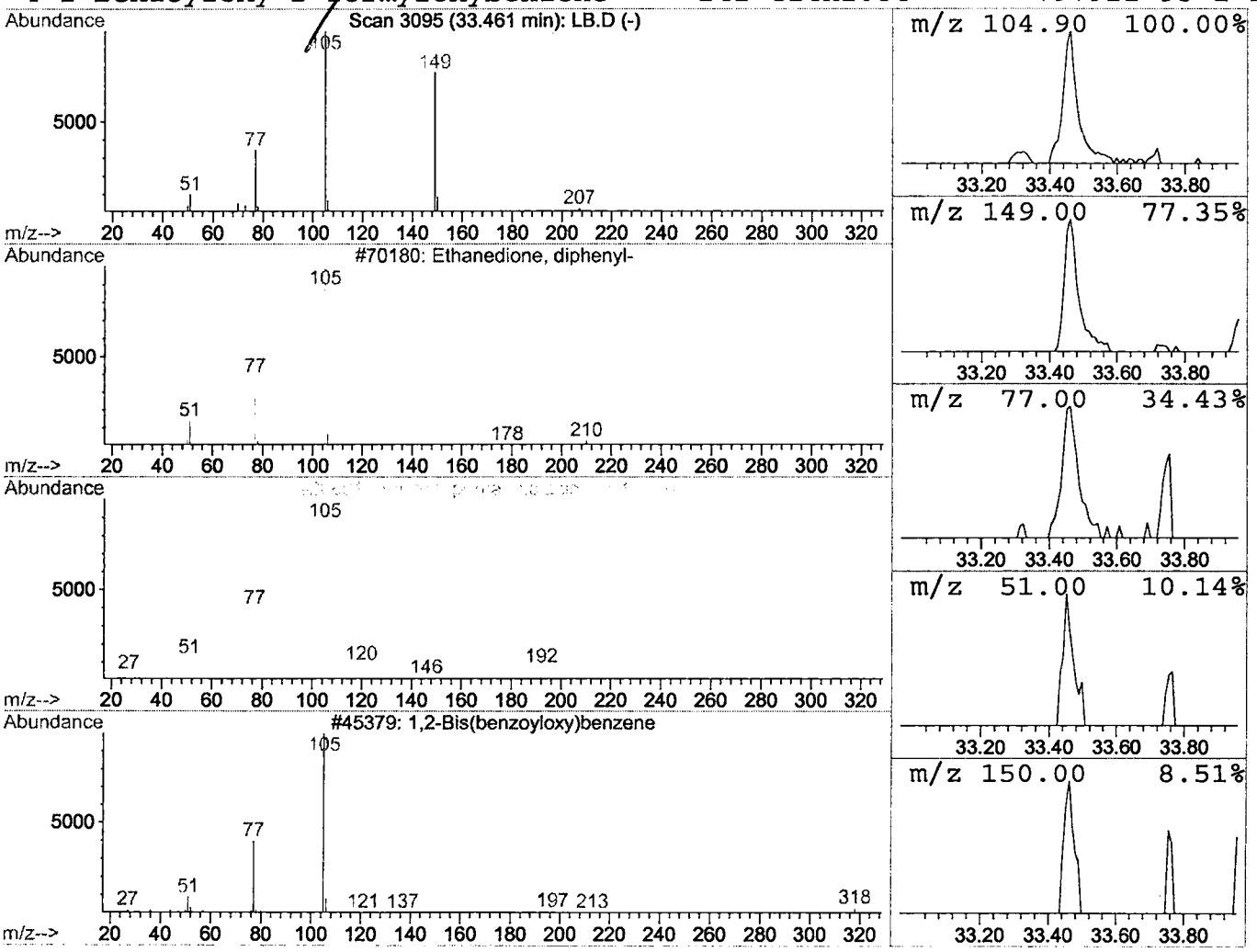
Quant Method : C:\HPCHEM\1\METHODS\GCMS0208.M (RTE Integrator)
Title : 209 625/TCLP calibration table
Library : C:\DATABASE\NBS75K.L

****

Peak Number 9 Ethanedione, diphenyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
33.46	0.22 ug/l	34987	Chrysene-d12	33.76

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Ethanedione, diphenyl-	210	C14H10O2	000134-81-6	10
2	Benzenepentanoic acid, .delta.-oxo-	192	C11H12O3	001501-05-9	10
3	1,2-Bis(benzoyloxy)benzene	318	C20H14O4	000643-94-7	10
4	1-Benzoyloxy-2-formyloxybenzene	242	C14H10O4	797922-93-1	10



Tentatively Identified Compound (LSC) summary

Operator ID: Date Acquired: 8 Mar 2002 3:29 am
Data File: C:\HPCHEM\1\DATA\MAR0702\LB.D
Name: gc/ms scan 2/5
Misc:
Method: C:\HPCHEM\1\METHODS\GCMS0208.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
Pentane, 3-methoxy-	8.20	0.5	ug/l	53325	ISTD01	12.30	2300970	20.0
Cyclopentasiloxane,	14.89	0.5	ug/l	71158	ISTD02	15.94	3046500	20.0
Cyclohexasiloxane, d	18.06	0.9	ug/l	132290	ISTD02	15.94	3046500	20.0
Trisiloxane, 1,1,1,5	20.88	0.7	ug/l	116503	ISTD03	21.25	3293410	20.0
Benzoic acid, 4-etho	21.65	0.3	ug/l	52141	ISTD03	21.25	3293410	20.0
3-Isopropoxy-1,1,1,7	23.41	0.5	ug/l	77217	ISTD03	21.25	3293410	20.0
1,1,1,5,7,7,7-Heptam	25.58	0.3	ug/l	51181	ISTD04	25.69	3379720	20.0
Trisiloxane, 1,1,1,5	27.52	0.2	ug/l	33925	ISTD04	25.69	3379720	20.0
Ethanedione, dipheny	33.46	0.2	ug/l	34987	ISTD05	33.76	3202650	20.0

column bleed

LB.D GCMS0208.M Wed Mar 13 15:04:09 2002