User: Fakhouri, Morgan
Westchester Dept. of Labs & Research
Date: 01/25/2002 Time: 14:22:37

Sample: AD30705
Analysis: $B1GCMS Blank for GCMS Scan
Units: ug
Started: 1/25/02 14:21 Ended: 1/25/02 14:21 Analyst: MF
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\DEC3101\LBE.D Vial: 40
Acq On : 2 Jan 2002 2:43 am Operator: 209 GALOS
Sample : gcms scan 12/19a Inst : GC/MS 209
Misc : Multiplr: 1.00
MS Integration Params: RTEINT.P
Quant Time: Jan 25 9:30 2002 Quant Results File: GCMS1126.RES

Quant Method : C:\HPCHEM\1\METHODS\GCMS1126.M (RTE Integrator)
Title : 209 625/TCLP calibration table
Last Update : Tue Jan 22 11:56:23 2002
Response via : Initial Calibration
DataAcq Meth : GCMS1126

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	11.67	152	1043144	20.00	ug/l	-0.03
10) Naphthalene-d8	15.28	136	3985818	20.00	ug/l	-0.03
17) Acenaphthene-d10	20.56	164	1984454	20.00	ug/l	-0.02
29) Phenanthrene-d10	24.98	188	3013816	20.00	ug/l	-0.01
38) Chrysene-d12	33.02	240	1925708	20.00	ug/l	-0.01
44) Perylene-d12	37.12	264	1152507	20.00	ug/l	0.00
System Monitoring Compounds						
37) 2,4-DDD	30.07	235	152327	4.60	ug/mL	0.00
Spiked Amount 5.000			Recovery = 92.00%			
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	14.20	82	614	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	15.33	128	1170	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	20.27	163	394	N.D.		
21) 2,6-Dinitrotoluene	0.00	165	0	N.D.		
22) Acenaphthylene	20.14	152	196	N.D.		
23) Acenaphthene	20.65	153	354	N.D.		
24) 2,4-Dinitrotoluene	21.34	165	197	N.D.		
25) Diethylphthalate	22.14	149	2361	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.		

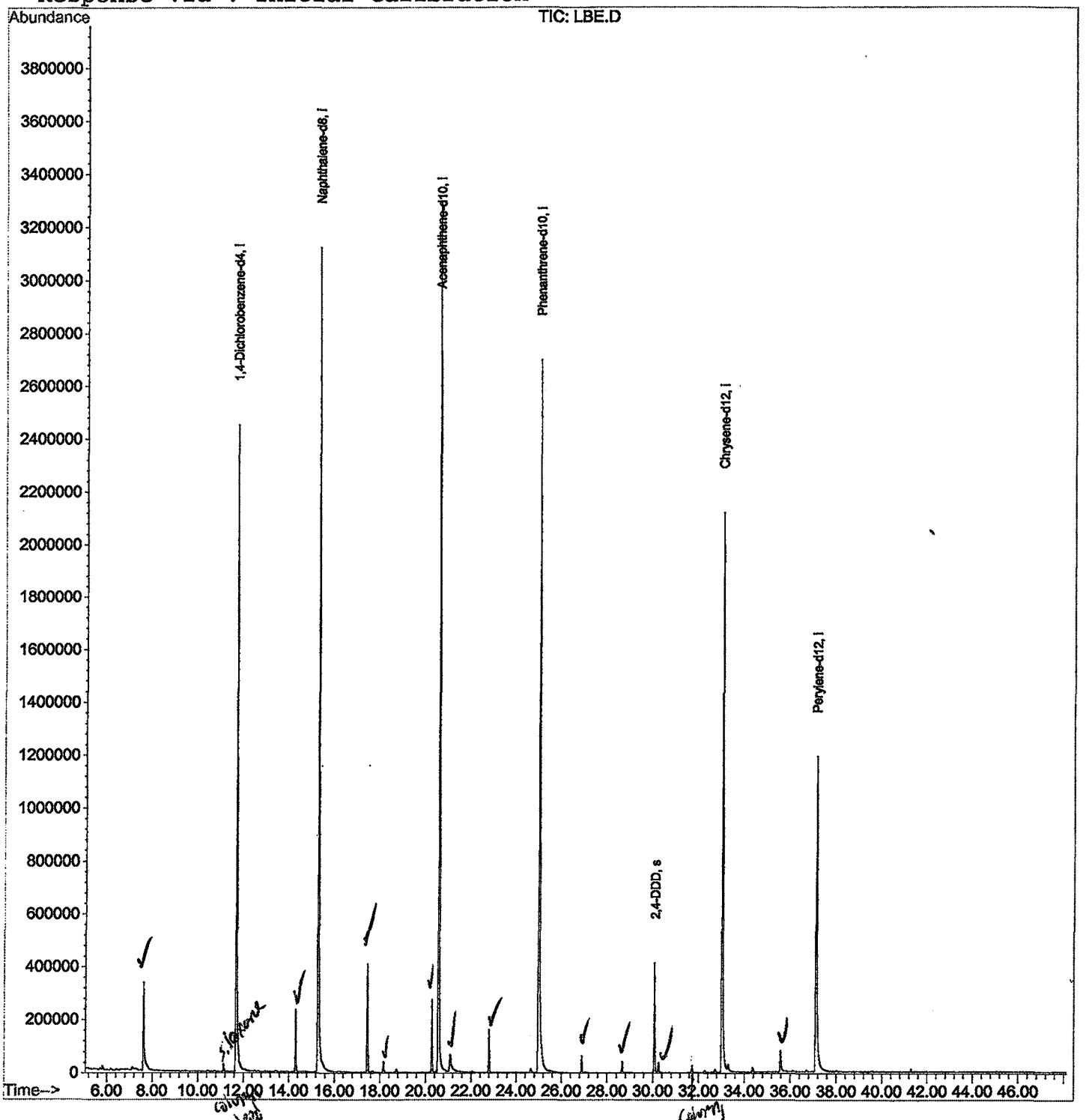
(#) = qualifier out of range (m) = manual integration

Quantitation Report

Data File : C:\HPCHEM\1\DATA\DEC3101\LBE.D
Acq On : 2 Jan 2002 2:43 am
Sample : gcms scan 12/19a
Misc :
MS Integration Params: RTEINT.P
Quant Time: Jan 25 9:30 2002

Vial: 40
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 : Tue Jan 22 11:56:23 2002
Response via : Initial Calibration



LSC Area Percent Report

Data File : C:\HPCHEM\1\DATA\DEC3101\LBE.D Vial: 40
Acq On : 2 Jan 2002 2:43 am Operator: 209 GALOS
Sample : gcms scan 12/19a 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

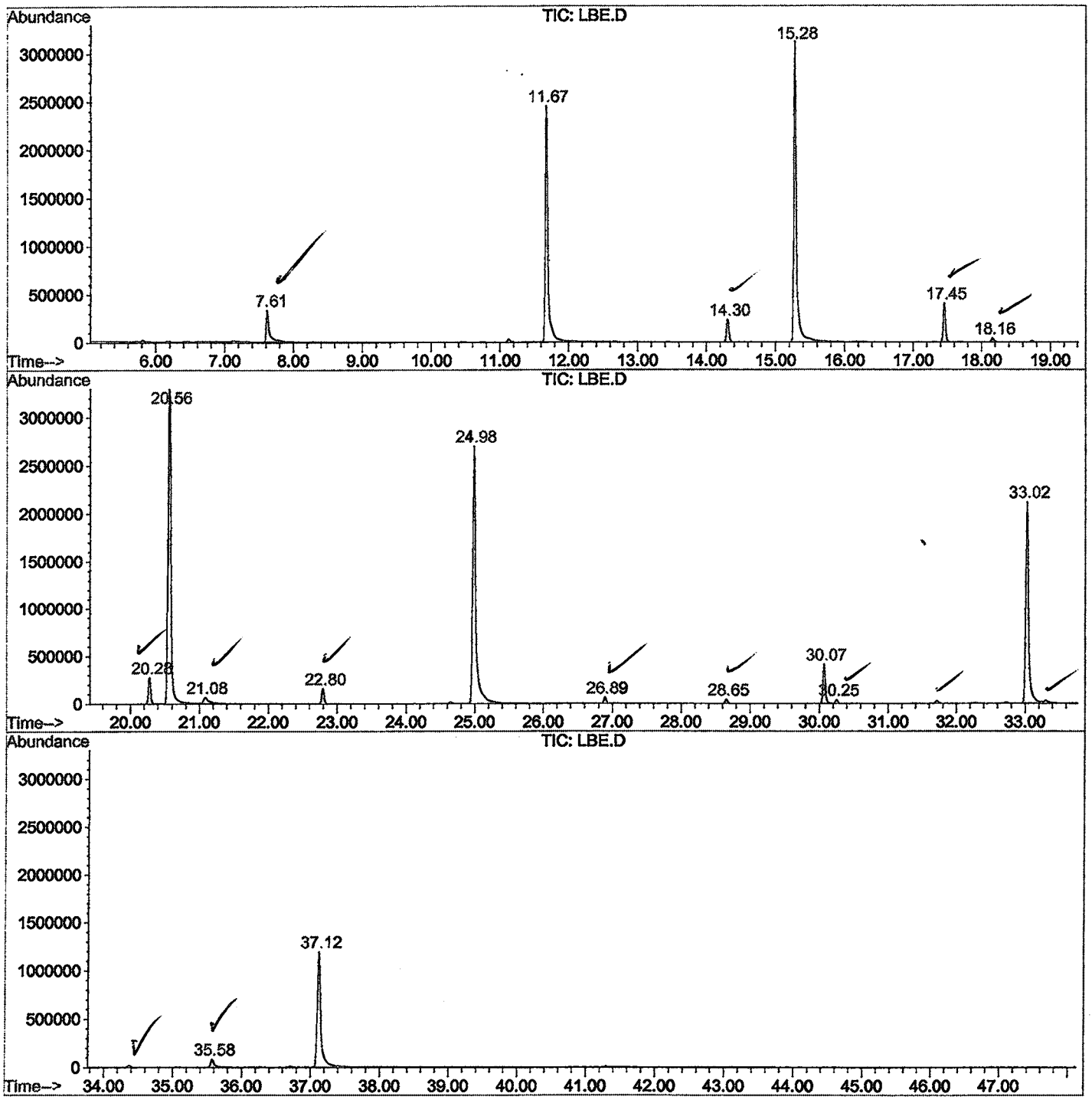
peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	peak area	peak % max.	% of total
1	7.612	276	280	314	rBV	331404	1067607	12.20%	2.219%
2	11.670	717	722	755	rBV	2449342	6540032	74.73%	13.592%
3	14.305	1003	1009	1023	rVB	234433	553224	6.32%	1.150%
4	15.278	1109	1115	1153	rBV	3121905	8208198	93.79%	17.059%
5	17.453	1346	1352	1360	rBV	407223	919355	10.51%	1.911%
6	18.160	1424	1429	1439	rBV	41587	103043	1.18%	0.214%
7	20.281	1654	1660	1668	rBV	275820	627578	7.17%	1.304%
8	20.556	1684	1690	1719	rBV	3297234	8751233	100.00%	18.188%
9	21.080	1740	1747	1767	rBV	62901	337199	3.85%	0.701%
10	22.796	1928	1934	1941	rBV	159180	368388	4.21%	0.766%
11	24.981	2164	2172	2208	rBV2	2698939	8645008	98.79%	17.967%
12	26.891	2374	2380	2387	rBV	63371	141727	1.62%	0.295%
13	28.653	2567	2572	2578	rBV	42590	94506	1.08%	0.196%
14	30.067	2719	2726	2740	rBV	413675	958820	10.96%	1.993%
15	30.251	2740	2746	2753	rVB	38613	89691	1.02%	0.186%
16	33.023	3041	3048	3069	rBV2	2117246	6173541	70.54%	12.830%
17	35.575	3316	3326	3341	rBV2	80345	290412	3.32%	0.604%
18	37.118	3486	3494	3522	rBV2	1189090	4246727	48.53%	8.826%

Sum of corrected areas: 48116289

LBE.D GCMS1126.M Fri Jan 25 12:07:06 2002

LSC Report - Integrated Chromatogram

File : C:\HPCHEM\1\DATA\DEC3101\LBE.D
Operator : 209 GALOS
Acquired : 2 Jan 2002 2:43 am using AcqMethod GCMS1126
Instrument : GC/MS 209
Sample Name: gcms scan 12/19a
Misc Info :
Vial Number: 40
Quant File :GCMS1126.RES (RTE Integrator)



LBE.D GCMS1126.M Fri Jan 25 12:07:12 2002

Library Search Compound Report

Data File : C:\HPCHEM\1\DATA\DEC3101\LBE.D
Acq On : 2 Jan 2002 2:43 am
Sample : gcms scan 12/19a
Misc :
MS Integration Params: LSCINT.P

Vial: 40
Operator: 209 GALOS
Inst : GC/MS 209
Multiplr: 1.00

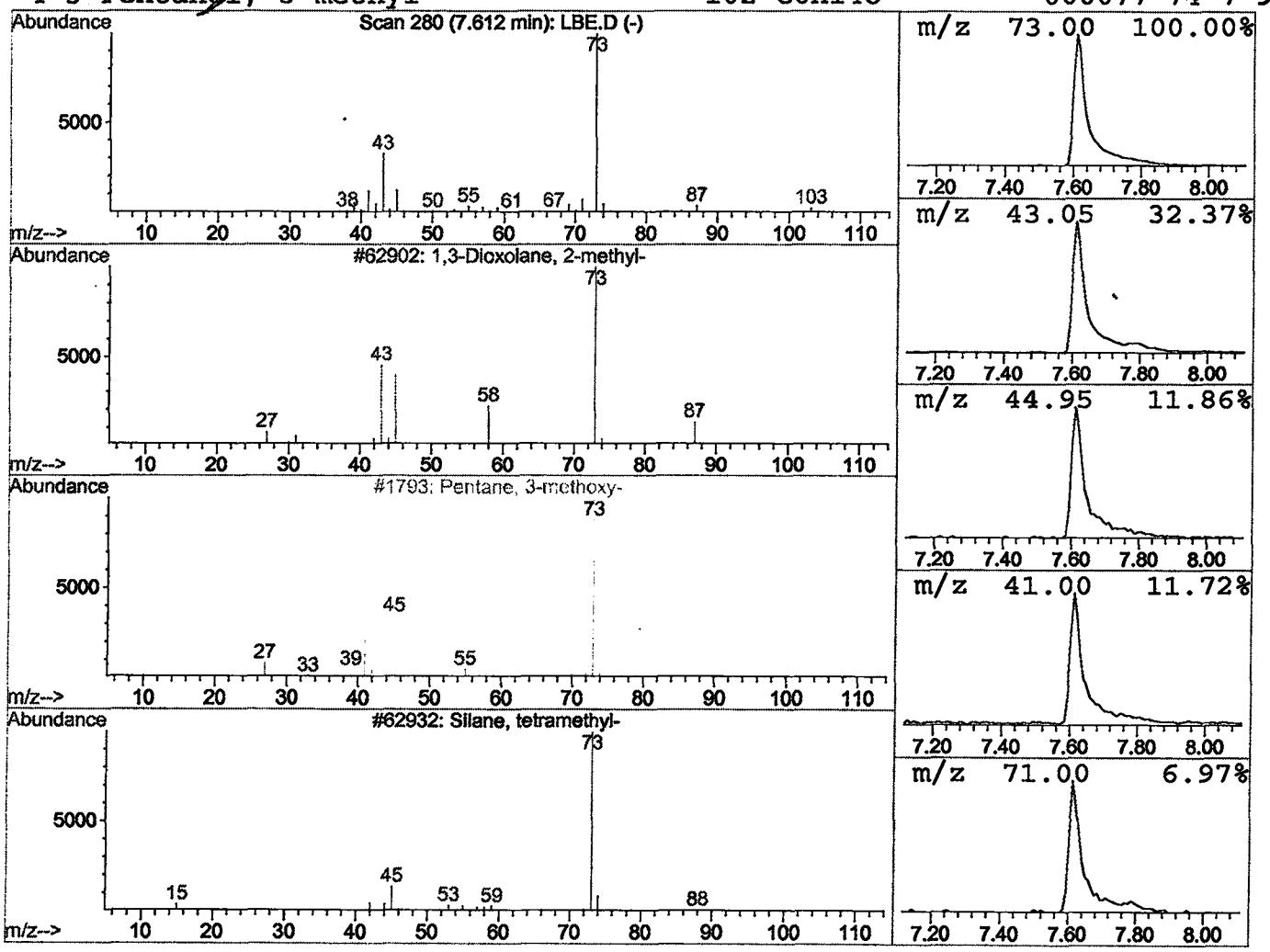
Quant Method : C:\HPCHEM\1\METHODS\GCMS1126.M (RTE Integrator)
Title : 209 625/TCLP calibration table
Library : C:\DATABASE\NBS75K.L

****

Peak Number 1 1,3-Dioxolane, 2-methyl- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.61	3.26 ug/l	1067610	1,4-Dichlorobenzene-d4	11.67

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,3-Dioxolane, 2-methyl-	88	C4H8O2	000497-26-7	9
2	Pentane, 3-methoxy-	102	C6H14O	036839-67-5	9
3	Silane, tetramethyl-	88	C4H12Si	000075-76-3	9
4	3-Pentanol, 3-methyl-	102	C6H14O	000077-74-7	9



Library Search Compound Report

Data File : C:\HPCHEM\1\DATA\DEC3101\LBE.D
Acq On : 2 Jan 2002 2:43 am
Sample : gcms scan 12/19a
Misc :
MS Integration Params: LSCINT.P

Vial: 40
Operator: 209 GALOS
Inst : GC/MS 209
Multiplr: 1.00

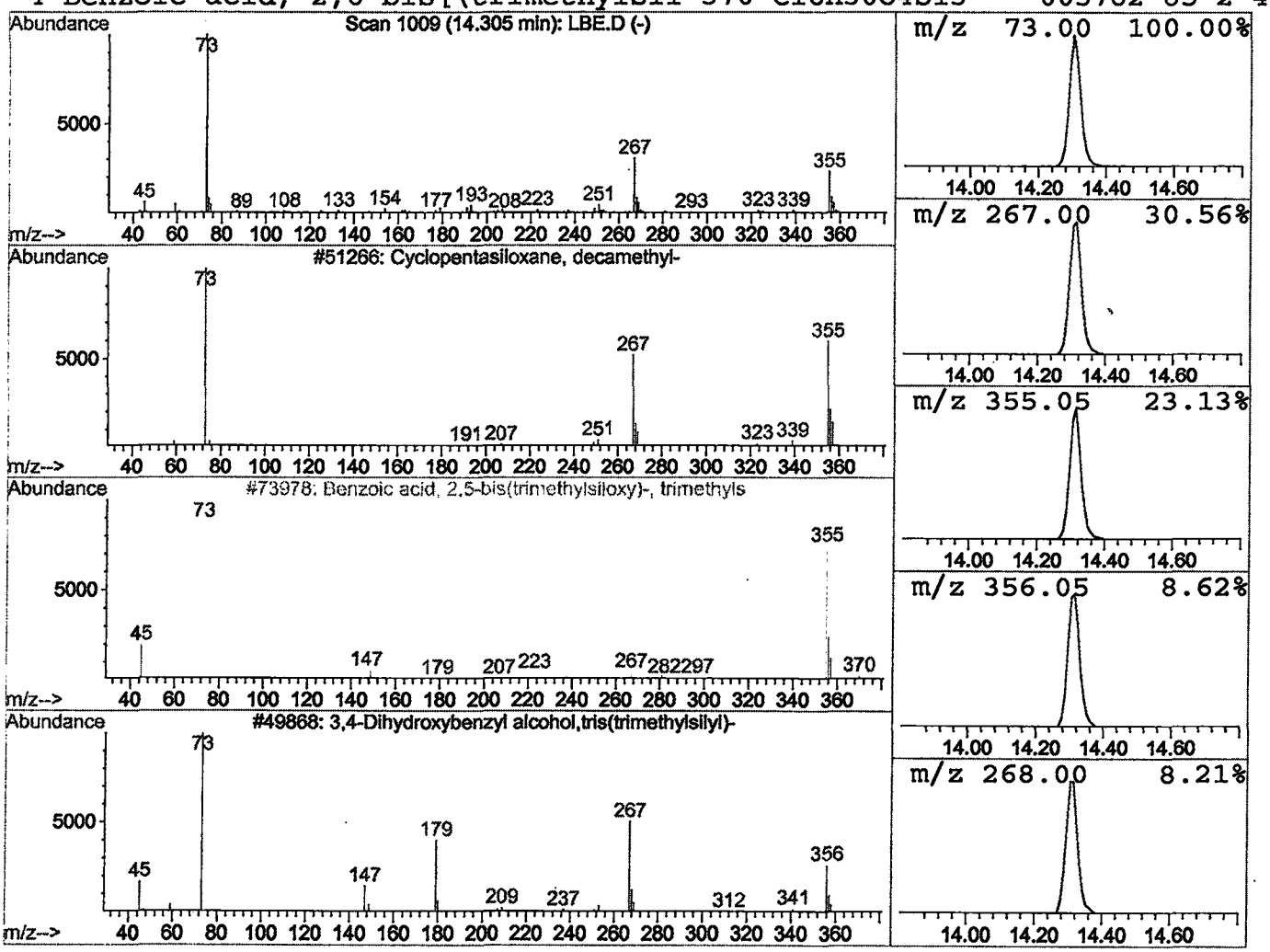
Quant Method : C:\HPCHEM\1\METHODS\GCMS1126.M (RTE Integrator)
Title : 209 625/TCLP calibration table
Library : C:\DATABASE\NBS75K.L

****

Peak Number 2 Cyclopentasiloxane, decamethyl Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.30	1.35 ug/l	553224	Naphthalene-d8	15.28

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclopentasiloxane, decamethyl-	370	C10H30O5Si5	000541-02-6	50
2	Benzoic acid, 2,5-bis(trimethylsilo	370	C16H30O4Si3	003618-20-0	47
3	3,4-Dihydroxybenzyl alcohol,tris(tr	356	C16H32O3Si3	068595-79-9	46
4	Benzoic acid, 2,6-bis[(trimethylsil	370	C16H30O4Si3	003782-85-2	40



Library Search Compound Report

Data File : C:\HPCHEM\1\DATA\DEC3101\LBE.D
Acq On : 2 Jan 2002 2:43 am
Sample : gcms scan 12/19a
Misc :
MS Integration Params: LSCINT.P

Vial: 40
Operator: 209 GALOS
Inst : GC/MS 209
Multiplr: 1.00

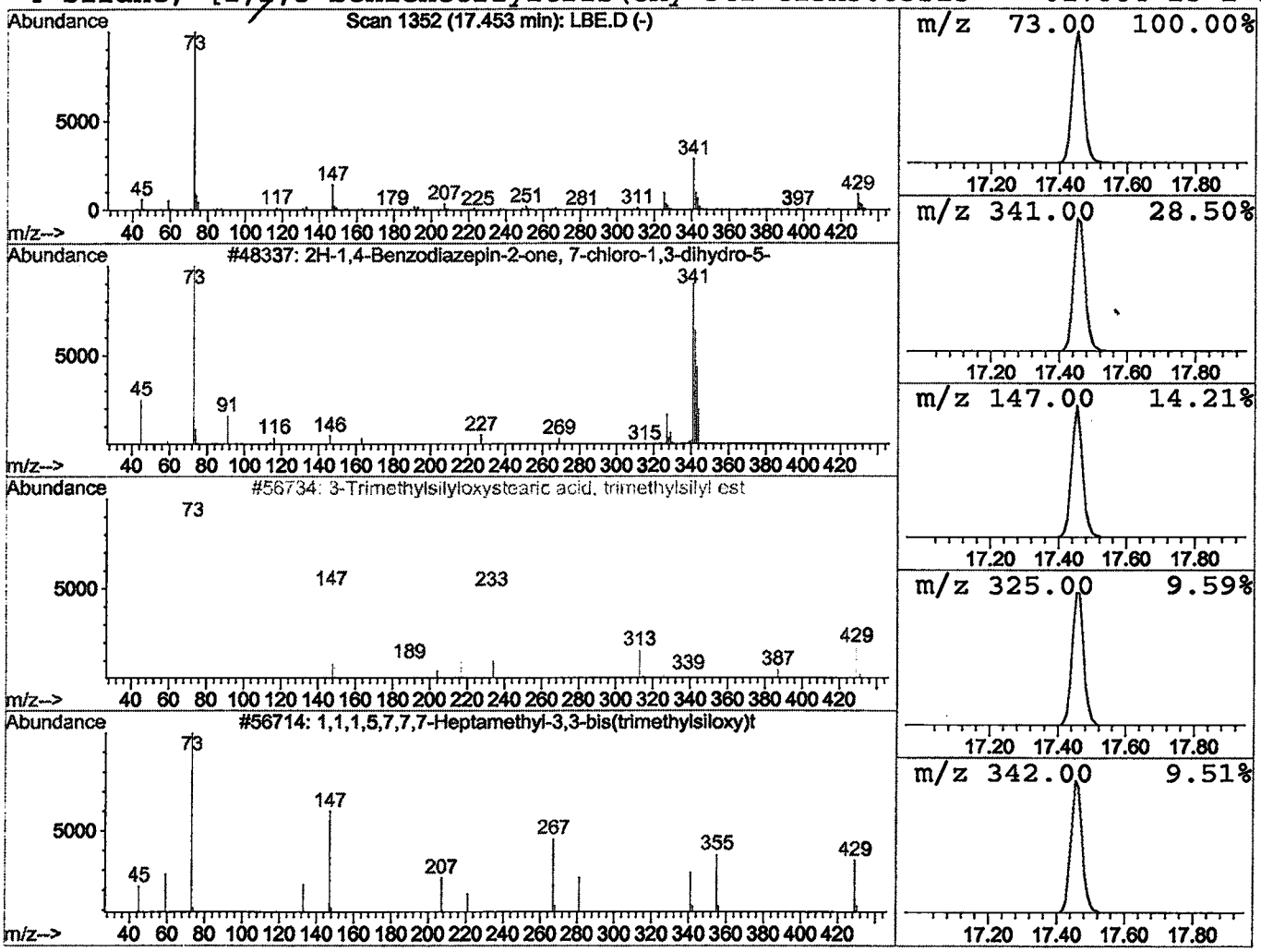
Quant Method : C:\HPCHEM\1\METHODS\GCMS1126.M (RTE Integrator)
Title : 209 625/TCLP calibration table
Library : C:\DATABASE\NBS75K.L

****

Peak Number 3 2H-1,4-Benzodiazepin-2-one, 7- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.45	2.24 ug/l	919355	Naphthalene-d8	15.28

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2H-1,4-Benzodiazepin-2-one, 7-chlor	342	C18H19ClN2OSi	055299-24-6	22
2	3-Trimethylsilyloxystearic acid, tr	444	C24H52O3Si2	000000-00-0	17
3	1,1,1,5,7,7,7-Heptamethyl-3,3-bis(t	444	C13H40O5Si6	038147-00-1	16
4	Silane, [1,2,3-benzenetriyltris(oxy	342	C15H30O3Si3	017864-23-2	10



Library Search Compound Report

Data File : C:\HPCHEM\1\DATA\DEC3101\LBE.D
Acq On : 2 Jan 2002 2:43 am
Sample : gcms scan 12/19a
Misc :
MS Integration Params: LSCINT.P

Vial: 40
Operator: 209 GALOS
Inst : GC/MS 209
Multiplr: 1.00

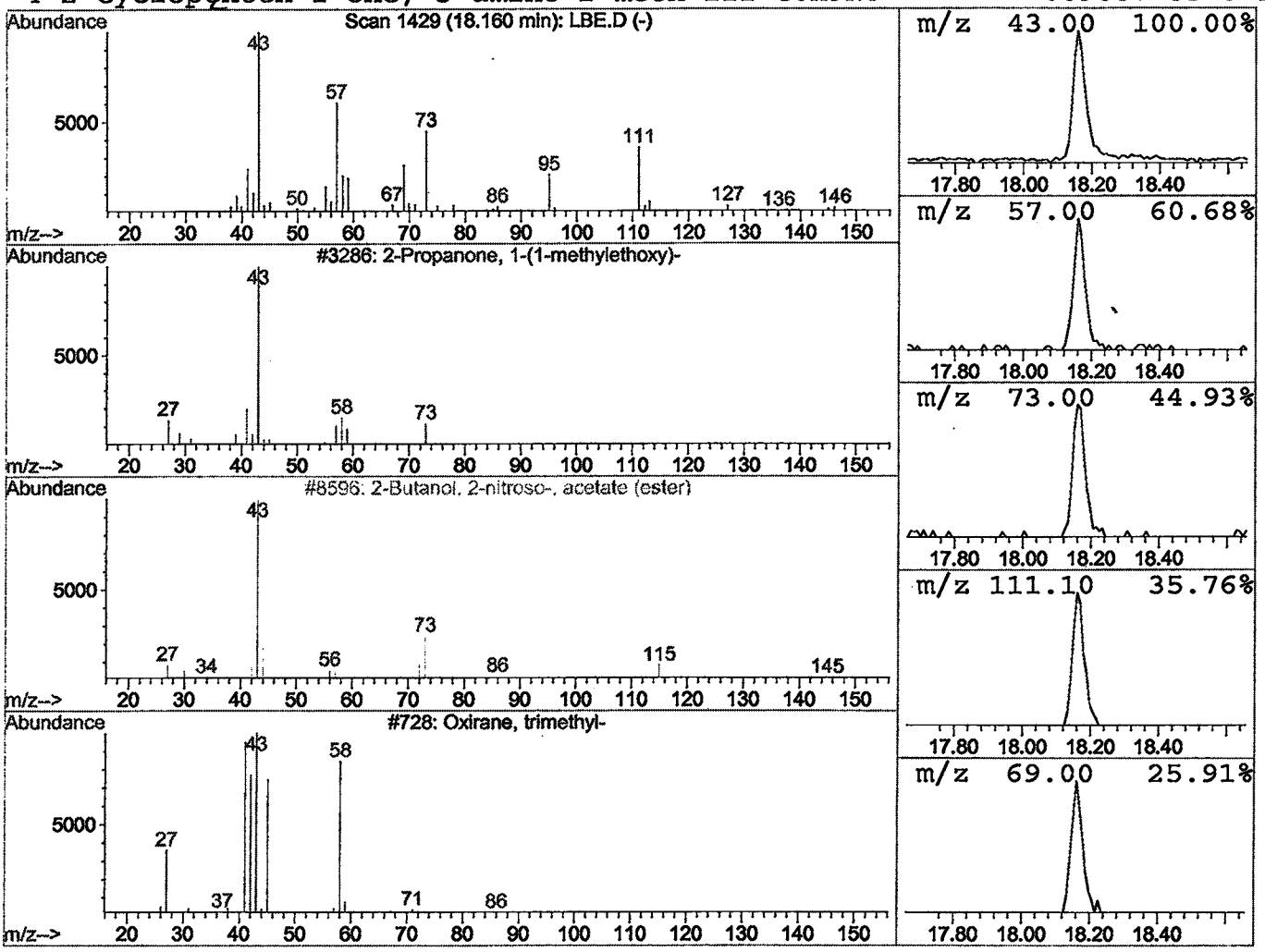
Quant Method : C:\HPCHEM\1\METHODS\GCMS1126.M (RTE Integrator)
Title : 209 625/TCLP calibration table
Library : C:\DATABASE\NBS75K.L

****

Peak Number 4 2-Propanone, 1-(1-methylethoxy Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.16	0.24 ug/l	103043	Acenaphthene-d10	20.56

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Propanone, 1-(1-methylethoxy)-	116	C6H12O2	042781-12-4	25
2	2-Butanol, 2-nitroso-, acetate (est	145	C6H11NO3	013880-90-5	9
3	Oxirane, trimethyl-	86	C5H10O	005076-19-7	9
4	2-Cyclopenten-1-one, 3-amino-2-meth	111	C6H9NO	069687-85-0	9



Library Search Compound Report

Data File : C:\HPCHEM\1\DATA\DEC3101\LBE.D
Acq On : 2 Jan 2002 2:43 am
Sample : gcms scan 12/19a
Misc :
MS Integration Params: LSCINT.P

Vial: 40
Operator: 209 GALOS
Inst : GC/MS 209
Multiplr: 1.00

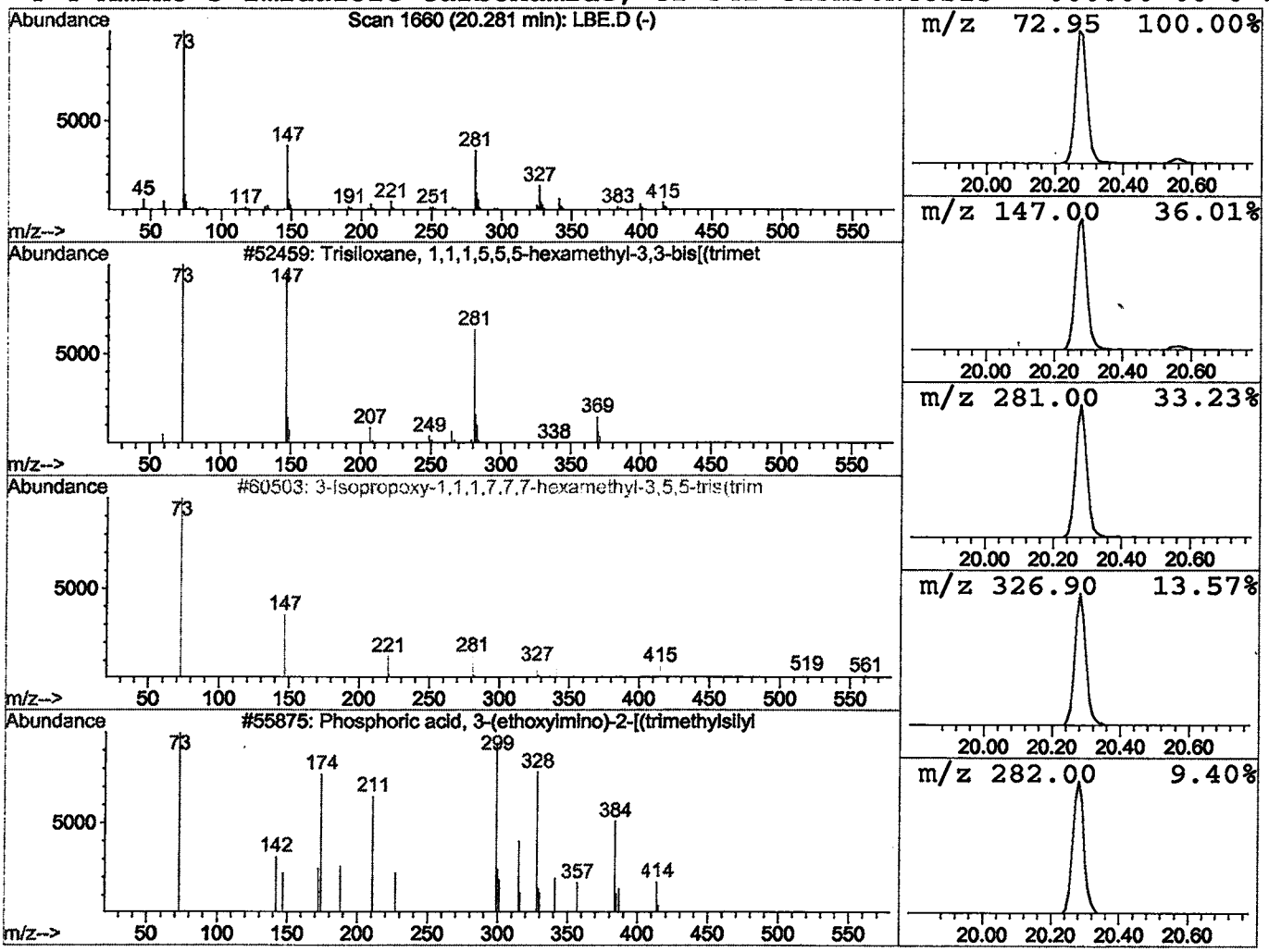
Quant Method : C:\HPCHEM\1\METHODS\GCMS1126.M (RTE Integrator)
Title : 209 625/TCLP calibration table
Library : C:\DATABASE\NBS75K.L

****

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.28	1.43 ug/l	627578	Acenaphthene-d10	20.56

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Trisiloxane, 1,1,1,5,5,5-hexamethyl	384	C12H36O4Si5	003555-47-3	42
2	3-Isopropoxy-1,1,1,7,7,7-hexamethyl	576	C18H52O7Si7	071579-69-6	38
3	Phosphoric acid, 3-(ethoxyimino)-2-	429	C14H36NO6PSi3	055334-96-8	14
4	4-Amino-5-imidazole carboxamide, tr	342	C13H30N4OSi3	000000-00-0	10



Library Search Compound Report

Data File : C:\HPCHEM\1\DATA\DEC3101\LBE.D
Acq On : 2 Jan 2002 2:43 am
Sample : gcms scan 12/19a
Misc :
MS Integration Params: LSCINT.P

Vial: 40
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

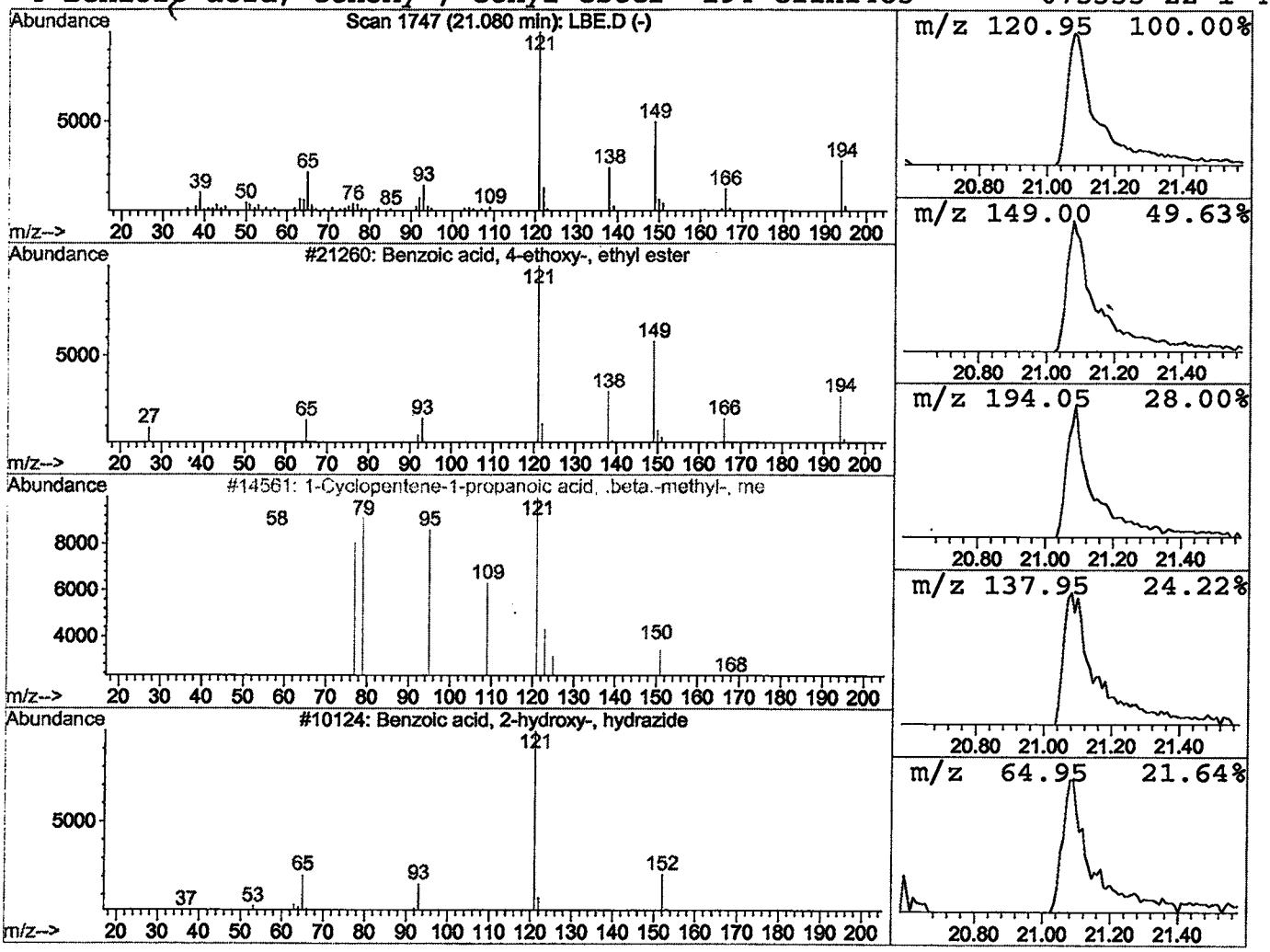
****

Peak Number 6 Benzoic acid, 4-ethoxy-, ethyl Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.08	0.77 ug/l	337199	Acenaphthene-d10	20.56

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	Benzoic acid, 2-hydroxy-, hydrazide	152	C7H8N2O2	000936-02-7	43
4	Benzoic acid, ethoxy-, ethyl ester	194	C11H14O3	075333-22-1	43

Column



Library Search Compound Report

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Acq On : 2 Jan 2002 2:43 am
Sample : gcms scan 12/19a
Misc :
MS Integration Params: LSCINT.P

Vial: 40
Operator: 209 GALOS
Inst : GC/MS 209
Multiplr: 1.00

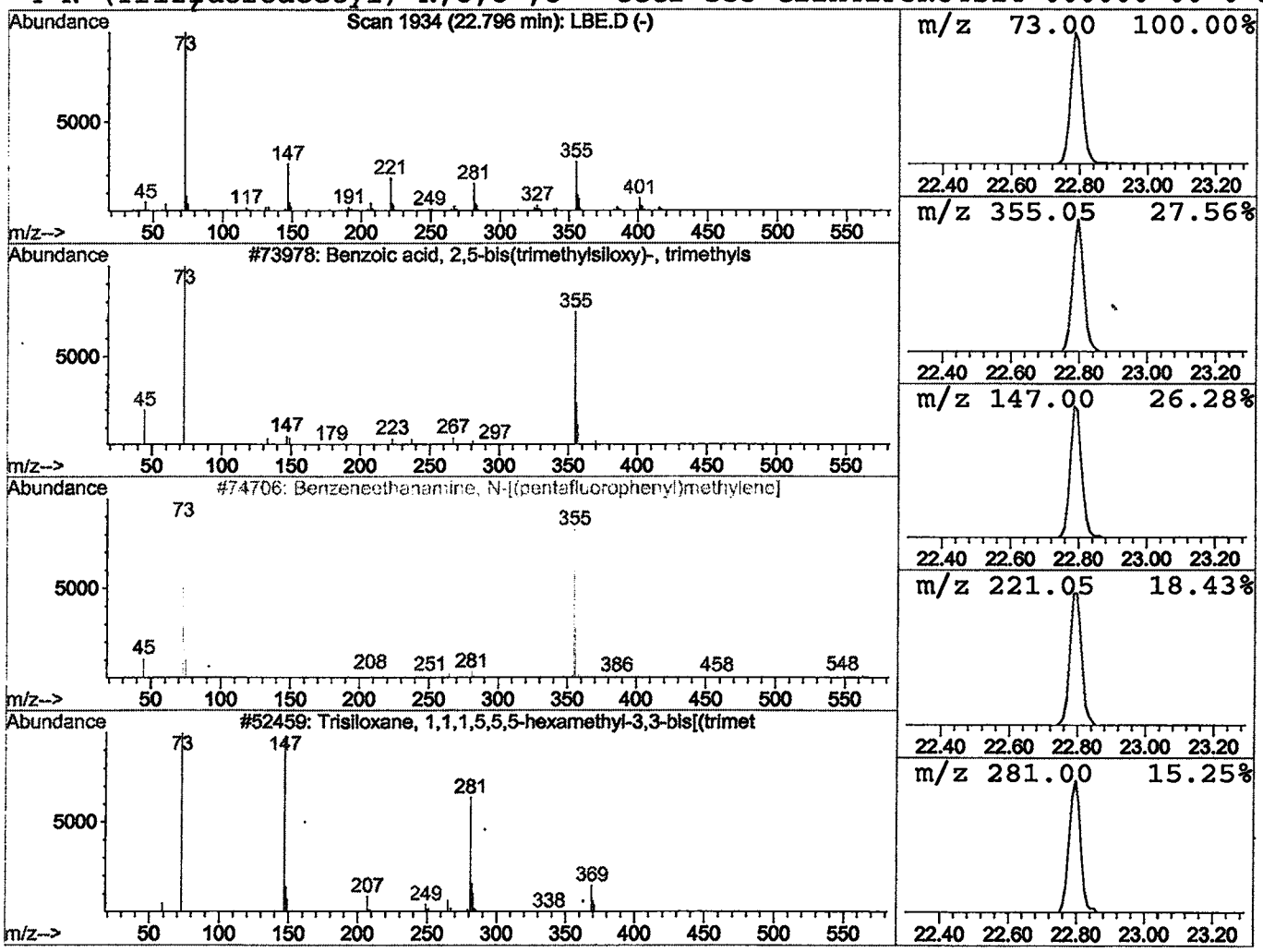
Quant Method : C:\HPCHEM\1\METHODS\GCMS1126.M (RTE Integrator)
Title : 209 625/TCLP calibration table
Library : C:\DATABASE\NBS75K.L

****

Peak Number 7 Benzoic acid, 2,5-bis(trimethy Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.80	0.85 ug/l	368388	Phenanthrene-d10	24.98

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzoic acid, 2,5-bis(trimethylsilo	370	C16H30O4Si3	003618-20-0	38
2	Benzeneethanamine, N-[(pentafluorop	563	C24H34F5NO3Si3	055429-13-5	37
3	Trisiloxane, 1,1,1,5,5,5-hexamethyl	384	C12H36O4Si5	003555-47-3	37
4	N-(Trifluoroacetyl)-N,O,O',O''-tetr	553	C22H42F3NO4Si4	000000-00-0	35



Library Search Compound Report

Data File : C:\HPCHEM\1\DATA\DEC3101\LBE.D
Acq On : 2 Jan 2002 2:43 am
Sample : gcms scan 12/19a
Misc :
MS Integration Params: LSCINT.P

Vial: 40
Operator: 209 GALOS
Inst : GC/MS 209
Multiplr: 1.00

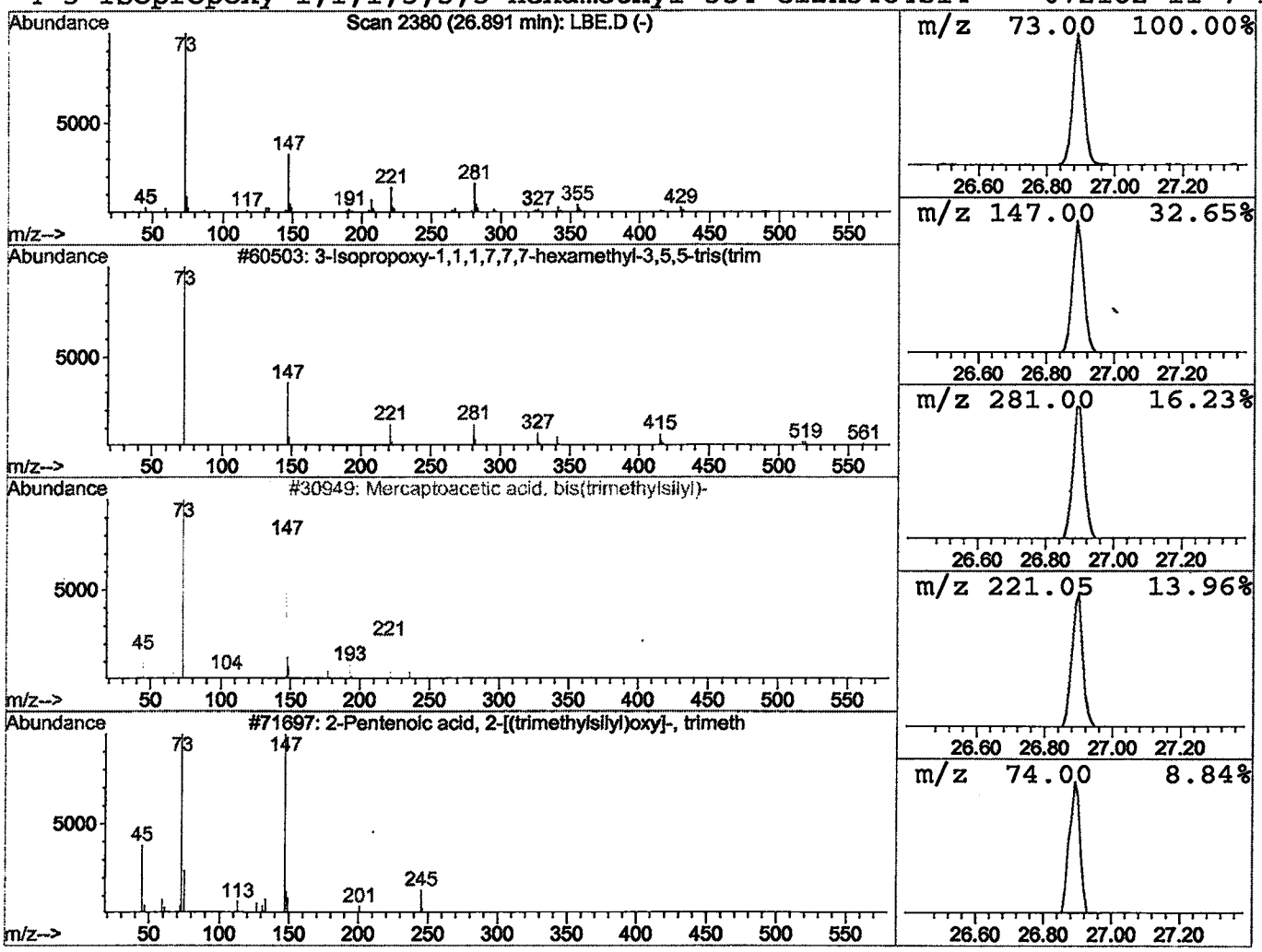
Quant Method : C:\HPCHEM\1\METHODS\GCMS1126.M (RTE Integrator)
Title : 209 625/TCLP calibration table
Library : C:\DATABASE\NBS75K.L

****

Peak Number 8 3-Isopropoxy-1,1,1,7,7,7-hexam Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
26.89	0.33 ug/l	141727	Phenanthrene-d10	24.98

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	3-Isopropoxy-1,1,1,7,7,7-hexamethyl	576	C18H52O7Si7	071579-69-6	28
2	Mercaptoacetic acid, bis(trimethyls	236	C8H20O2SSi2	006398-62-5	23
3	2-Pentenoic acid, 2-[(trimethylsily	260	C11H24O3Si2	055045-17-5	23
4	3-Isopropoxy-1,1,1,5,5,5-hexamethyl	354	C12H34O4Si4	072182-11-7	14



Library Search Compound Report

Data File : C:\HPCHEM\1\DATA\DEC3101\LBE.D
Acq On : 2 Jan 2002 2:43 am
Sample : gcms scan 12/19a
Misc :
MS Integration Params: LSCINT.P

Vial: 40
Operator: 209 GALOS
Inst : GC/MS 209
Multiplr: 1.00

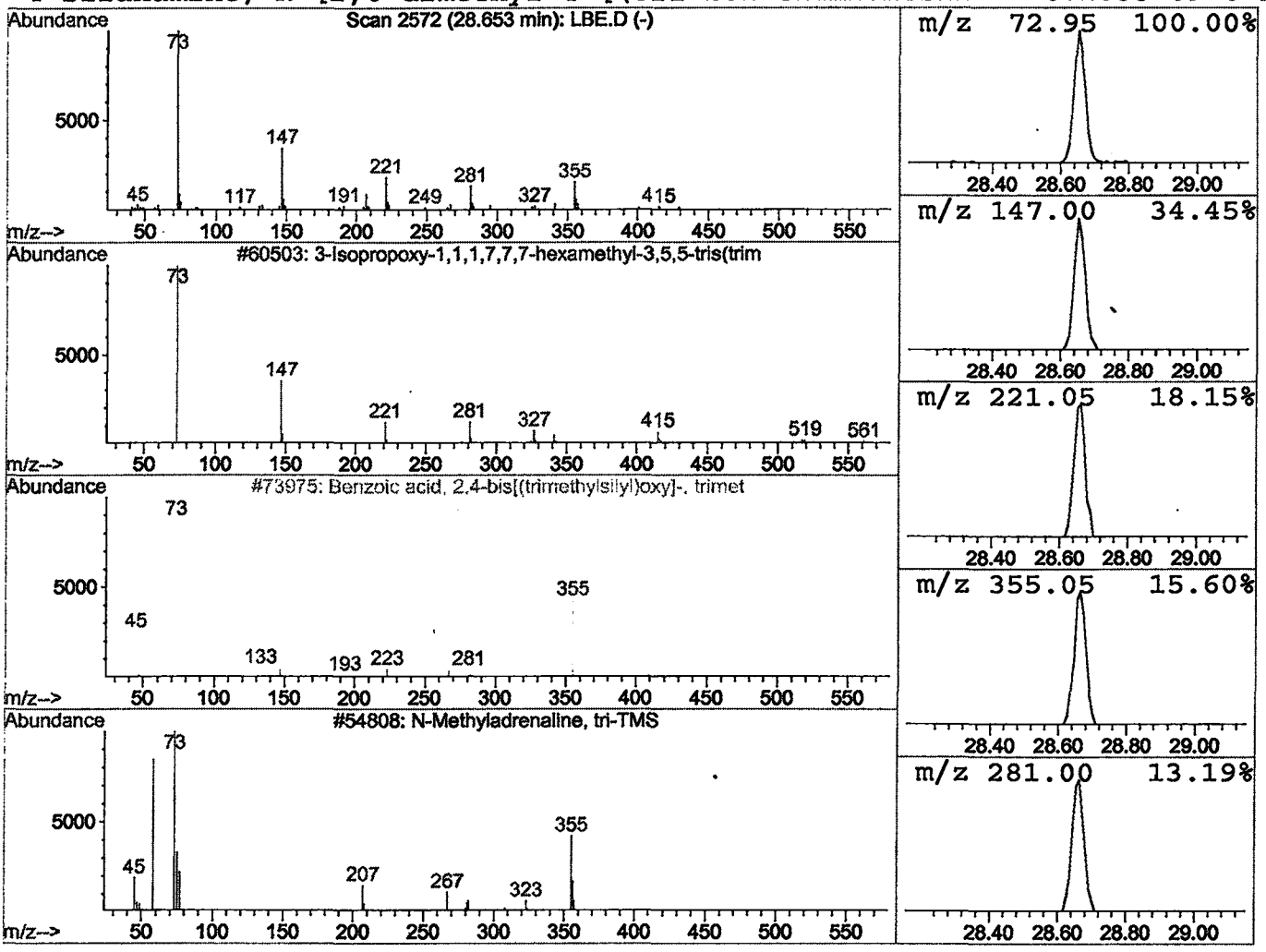
Quant Method : C:\HPCHEM\1\METHODS\GCMS1126.M (RTE Integrator)
Title : 209 625/TCLP calibration table
Library : C:\DATABASE\NBS75K.L

****

Peak Number 9 3-Isopropoxy-1,1,1,7,7,7-hexam Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
28.65	0.22 ug/l	94506	Phenanthrene-d10	24.98

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	3-Isopropoxy-1,1,1,7,7,7-hexamethyl	576	C18H52O7Si7	071579-69-6	42
2	Benzoic acid, 2,4-bis[(trimethylsil	370	C16H30O4Si3	010586-16-0	25
3	N-Methyladrenaline, tri-TMS	413	C19H39NO3Si3	000000-00-0	23
4	Silanamine, N-[2,6-dimethyl-4-[(tri	281	C14H27NOSi2	072088-09-6	9



Library Search Compound Report

Data File : C:\HPCHEM\1\DATA\DEC3101\LBE.D
Acq On : 2 Jan 2002 2:43 am
Sample : gcms scan 12/19a
Misc :
MS Integration Params: LSCINT.P

Vial: 40
Operator: 209 GALOS
Inst : GC/MS 209
Multiplr: 1.00

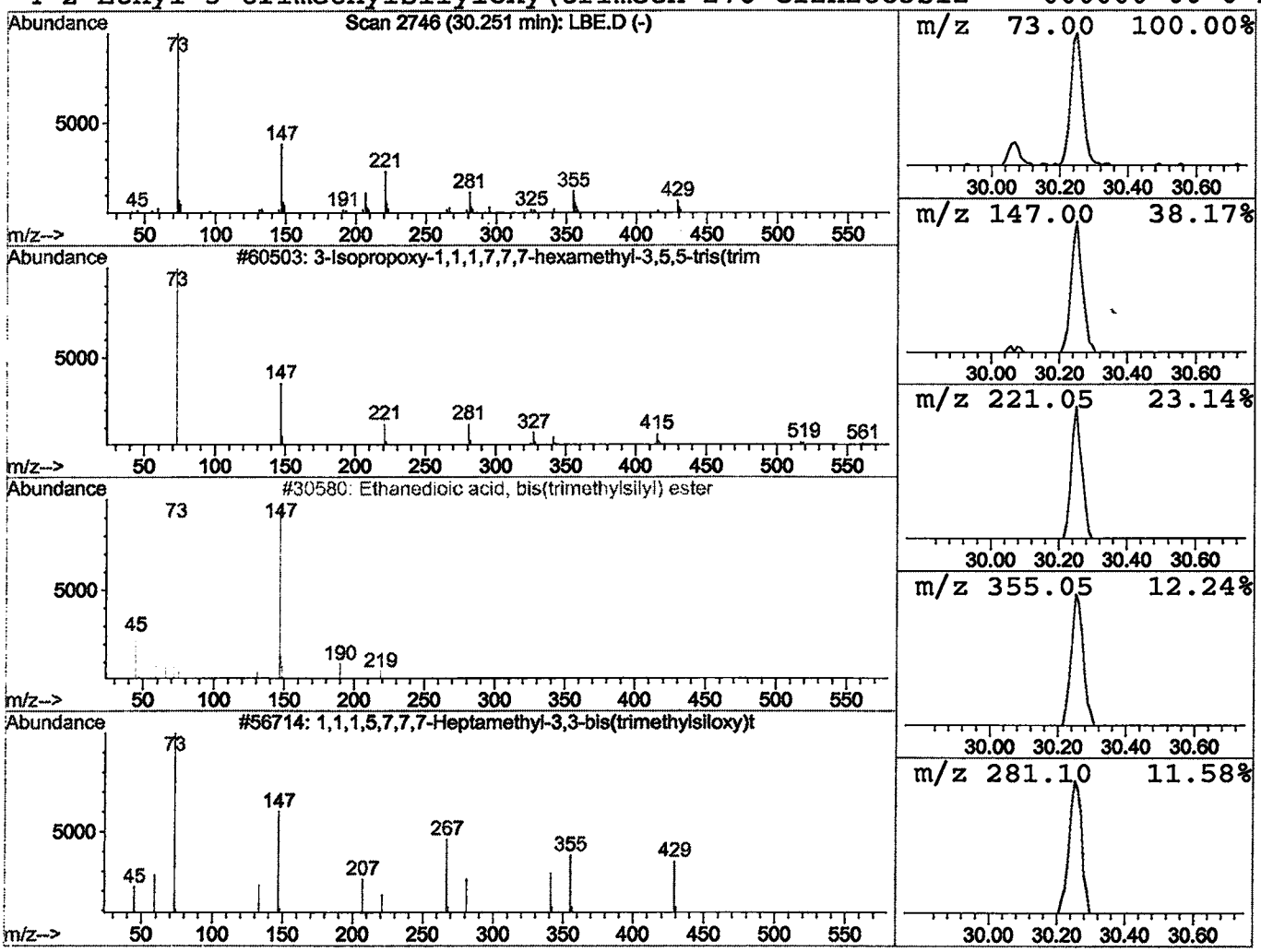
Quant Method : C:\HPCHEM\1\METHODS\GCMS1126.M (RTE Integrator)
Title : 209 625/TCLP calibration table
Library : C:\DATABASE\NBS75K.L

****

Peak Number 10 3-Isopropoxy-1,1,1,7,7,7-hexam Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
30.25	0.29 ug/l	89691	Chrysene-d12	33.02

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	3-Isopropoxy-1,1,1,7,7,7-hexamethyl	576	C18H52O7Si7	071579-69-6	40
2	Ethanedioic acid, bis(trimethylsily	234	C8H18O4Si2	018294-04-7	38
3	1,1,1,5,7,7,7-Heptamethyl-3,3-bis(t	444	C13H40O5Si6	038147-00-1	25
4	2-Ethyl-3-trimethylsilyloxy(trimeth	276	C12H28O3Si2	000000-00-0	23



Library Search Compound Report

Data File : C:\HPCHEM\1\DATA\DEC3101\LBE.D
Acq On : 2 Jan 2002 2:43 am
Sample : gcms scan 12/19a
Misc :
MS Integration Params: LSCINT.P

Vial: 40
Operator: 209 GALOS
Inst : GC/MS 209
Multiplr: 1.00

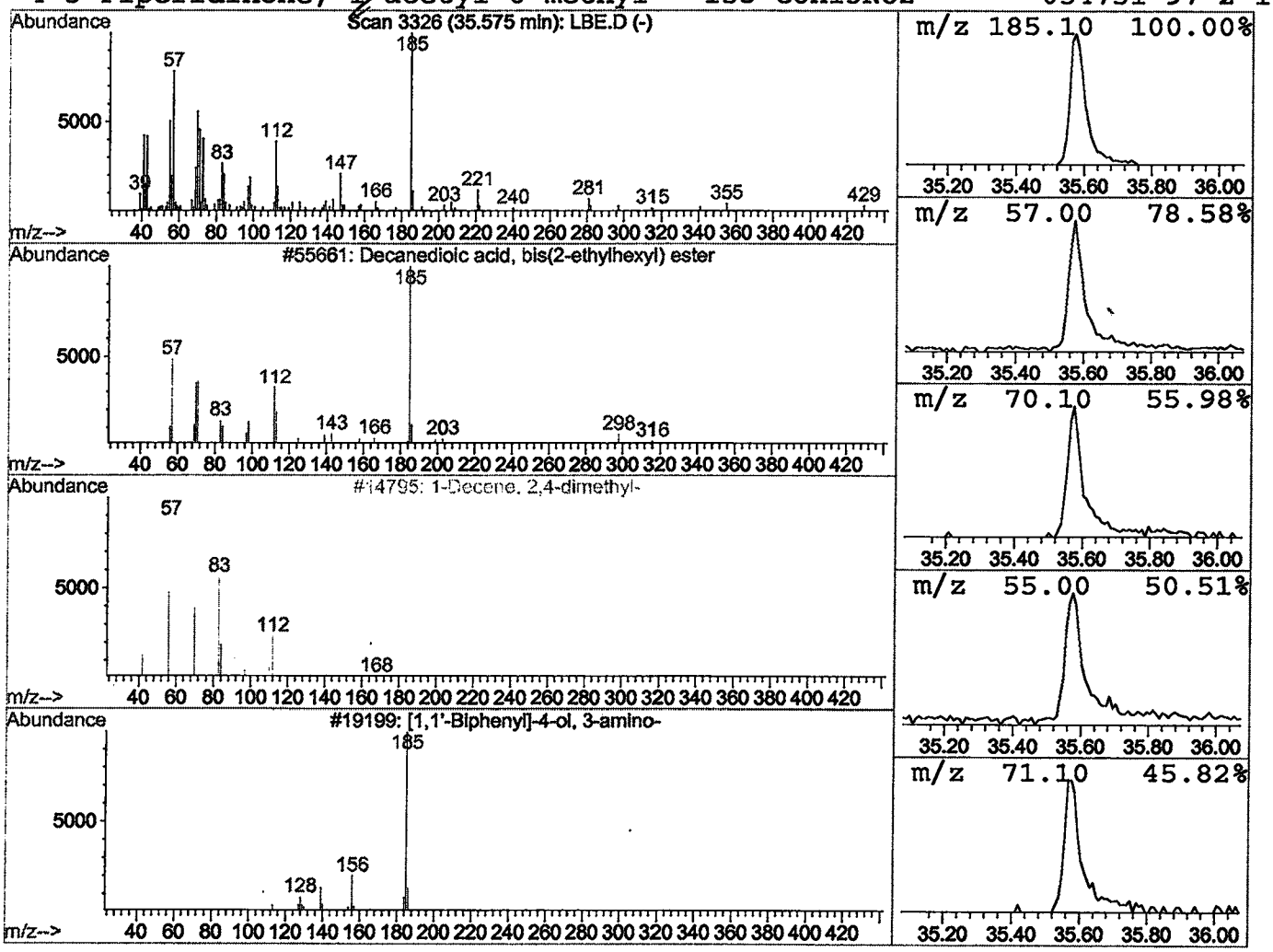
Quant Method : C:\HPCHEM\1\METHODS\GCMS1126.M (RTE Integrator)
Title : 209 625/TCLP calibration table
Library : C:\DATABASE\NBS75K.L

****

Peak Number 11 Decanedioic acid, bis(2-ethylh Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
35.58	1.37 ug/l	290412	Perylene-d12	37.12

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Decanedioic acid, bis(2-ethylhexyl)	426	C26H50O4	000122-62-3	59
2	1-Decene, 2,4-dimethyl-	168	C12H24	055170-80-4	25
3	[1,1'-Biphenyl]-4-ol, 3-amino-	185	C12H11NO	001134-36-7	22
4	3-Piperidinone, 1-acetyl-6-methyl-	155	C8H13NO2	054751-97-2	14



Tentatively Identified Compound (LSC) summary

Operator ID: 209 GALOS Date Acquired: 2 Jan 2002 2:43 am
Data File: C:\HPCHEM\1\DATA\DEC3101\LBE.D
Name: gcms scan 12/19a
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
1,3-Dioxolane, 2-met	7.61	3.3	ug/l	1067610	ISTD01	11.67	6540030	20.0
Cyclopentasiloxane,	14.30	1.3	ug/l	553224	ISTD02	15.28	8208200	20.0
2H-1,4-Benzodiazepin	17.45	2.2	ug/l	919355	ISTD02	15.28	8208200	20.0
2-Propanone, 1-(1-me	18.16	0.2	ug/l	103043	ISTD03	20.56	8751230	20.0
Trisiloxane, 1,1,1,5	20.28	1.4	ug/l	627578	ISTD03	20.56	8751230	20.0
Benzoic acid, 4-etho	21.08	0.8	ug/l	337199	ISTD03	20.56	8751230	20.0
Benzoic acid, 2,5-bi	22.80	0.9	ug/l	368388	ISTD04	24.98	8645010	20.0
3-Isopropoxy-1,1,1,7	26.89	0.3	ug/l	141727	ISTD04	24.98	8645010	20.0
3-Isopropoxy-1,1,1,7	28.65	0.2	ug/l	94506	ISTD04	24.98	8645010	20.0
3-Isopropoxy-1,1,1,7	30.25	0.3	ug/l	89691	ISTD05	33.02	6173540	20.0
Decanedioic acid, bi	35.58	1.4	ug/l	290412	ISTD06	37.12	4246730	20.0

LBE.D GCMS1126.M Fri Jan 25 12:08:08 2002