User: Vinci, David L
Westchester Dept. of Labs & Research
Date: 01/22/2002 Time: 12:03:24

Sample: AD30608
Analysis: $GCMS GCMS Scan For Unknowns
Units: ug
Started: 1/22/02 12:03 Ended: 1/22/02 12:03 Analyst: DLV
Page: 1

	Component Name	Viol	Result	Qualifier	MDL
1	Other unknown compounds		.		
2	Surr_2,4-DDD		90		5.0

Comments for Sample AD30608 - Analysis $GCMS

Westchester Dept. of Labs & Research
User: Vinci, David L
Date: 01-22-2002 Time: 12:03:47

The GCMS scan, from 35 to 550 amu, does not reveal the presence of unusual compounds.

Quantitation Report (QT Reviewed)

Data File : C:\HPCHEM\1\DATA\DEC2701\30608.D Vial: 32
Acq On : 28 Dec 2001 8:16 pm Operator: 209 GALOS
Sample : GC/MS SCAN 12/17 Inst : GC/MS 209
Misc : Multiplr: 1.00
MS Integration Params: GOOFY.P
Quant Time: Dec 31 14:10 2001 Quant Results File: GCMS1126.RES

Quant 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
DataAcq Meth : GCMS1126

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	11.68	152	1021716	20.00	ug/l	-0.02
10) Naphthalene-d8	15.28	136	3965981	20.00	ug/l	-0.02
17) Acenaphthene-d10	20.56	164	1978494	20.00	ug/l	-0.01
29) Phenanthrene-d10	24.99	188	2901181m	20.00	ug/l	0.00
38) Chrysene-d12	33.02	240	1801278	20.00	ug/l	-0.01
44) Perylene-d12	37.11	264	1090104	20.00	ug/l	0.00
System Monitoring Compounds						
37) 2,4-DDD	30.06	235	143965	4.51	ug/mL	0.00
Spiked Amount 5.000			Recovery = 90.20%			
Target Compounds						Qvalue
2) N-nitroso-dimethyl amine	5.17	74	1585	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	12.94	77	205	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	15.34	128	1552	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	20.28	163	324	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	21.24	165	182	N.D.		
25) Diethylphthalate	22.12	149	901	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\DEC2701\30608.D Vial: 32
Acq On : 28 Dec 2001 8:16 pm Operator: 209 GALOS
Sample : GC/MS SCAN 12/17 Inst : GC/MS 209
Misc : Multiplr: 1.00
MS Integration Params: GOOFY.P
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Quant 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
DataAcq Meth : GCMS1126

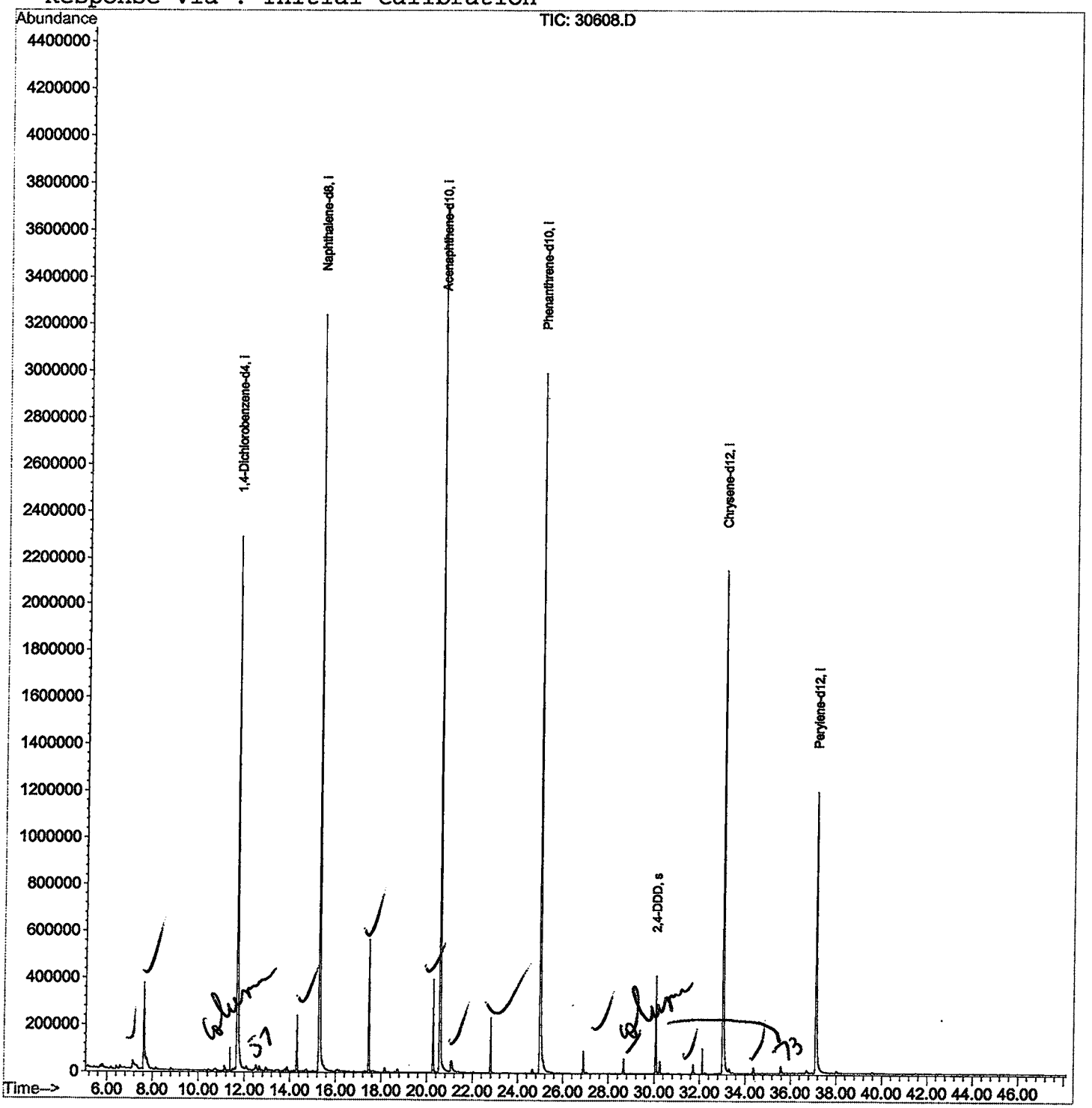
Compound	R.T.	QIon	Response	Conc Unit	Qvalue
30) N-Nitrosodiphenylamine	0.00	168	0	N.D.	
31) 4-Bromophenyl-phenylether	0.00	248	0	N.D.	
32) Hexachlorobenzene	0.00	284	0	N.D.	
33) Phenanthrene	25.04	178	1121	N.D.	
34) Anthracene	25.04	178	1121	N.D.	
35) Di-n-butylphthalate	27.03	149	8545	N.D.	
36) Fluoranthene	0.00	202	0	N.D.	
39) Pyrene	0.00	202	0	N.D.	
40) Butylbenzylphthalate	31.72	149	962	N.D.	
41) Benzo(a)anthracene	33.02	228	4996	N.D.	
42) Bis(2-ethylhexyl)phthalate	33.30	149	10683	N.D.	
43) Chrysene	33.02	228	4996	N.D.	
45) Di-n-Octylphthalate	35.04	149	1255	N.D.	
46) Benzo(b)fluoranthene	36.14	252	1248	N.D.	
47) Benzo(k)fluoranthene	36.10	252	492	N.D.	
48) Benzo(a)pyrene	36.96	252	343	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\DEC2701\30608.D Vial: 32
Acq On : 28 Dec 2001 8:16 pm Operator: 209 GALOS
Sample : GC/MS SCAN 12/17 Inst : GC/MS 209
Misc : Multiplr: 1.00
MS Integration Params: GOOFY.P
Quant Time: Dec 31 14:10 2001 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 : M:\INSTRU~1\209\DATA\DEC2701\30608.D Vial: 32
Acq On : 28 Dec 2001 8:16 pm Operator: 209 GALOS
Sample : GC/MS SCAN 12/17 Inst : GC/MS 209
Misc : Multiplr: 1.00
MS Integration Params: LSCINT.P

Method : C:\HPCHEM\1\METHODS\GCMS0103.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

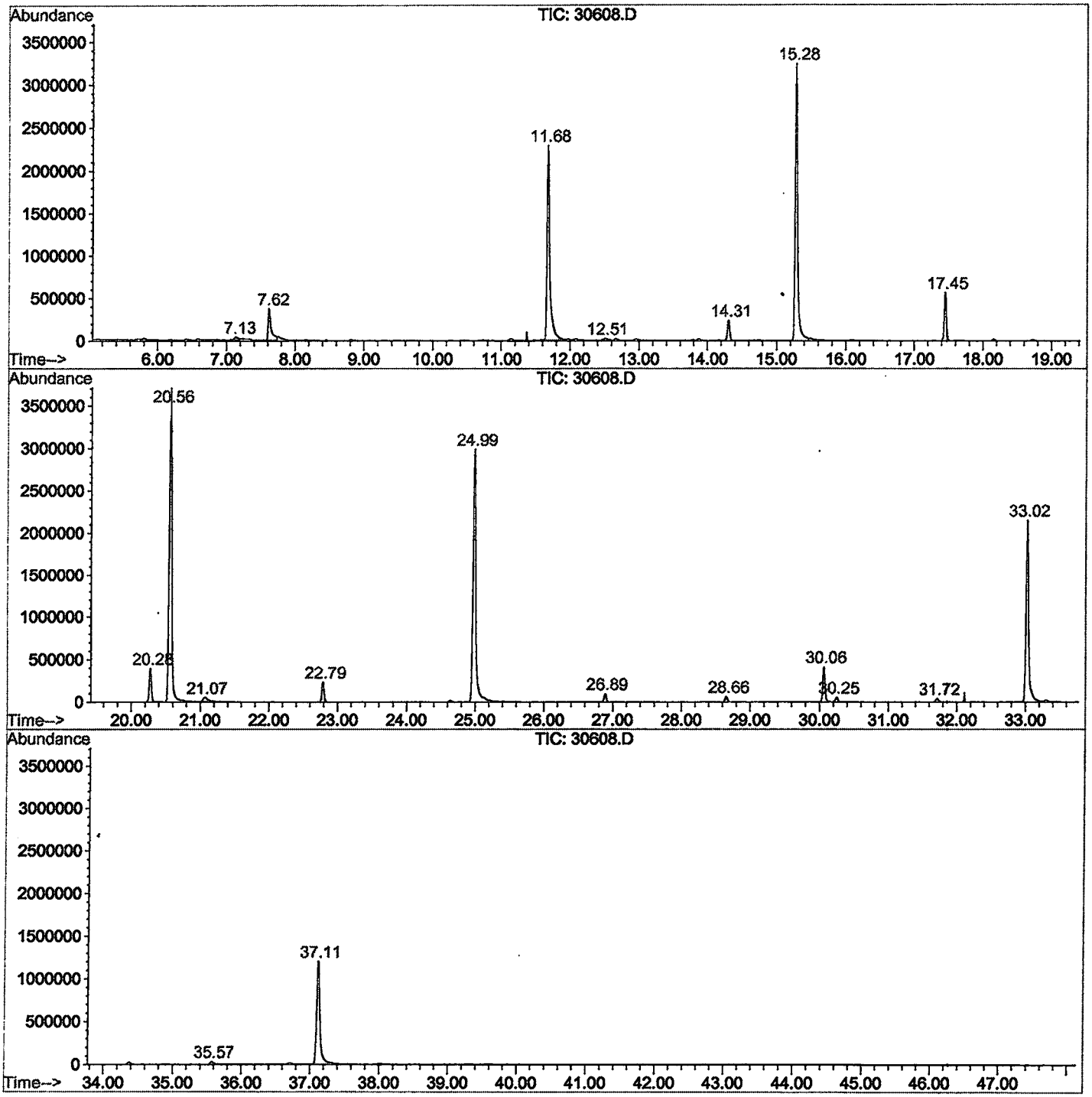
peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	peak area	peak % max.	% of total
1	7.131	222	227	234	rBV3	33936	130123	1.51%	0.274%
2	7.618	276	280	293	rBV	367941	1116872	13.00%	2.356%
3	11.676	717	722	752	rBV	2282452	6351977	73.95%	13.397%
4	12.511	805	813	824	rVB	23266	97254	1.13%	0.205%
5	14.310	999	1009	1017	rBV	235656	518333	6.03%	1.093%
6	15.284	1109	1115	1134	rBV	3236629	7936867	92.40%	16.740%
7	17.450	1345	1351	1362	rBV	561530	1193222	13.89%	2.517%
8	20.278	1652	1659	1667	rBV	394688	839991	9.78%	1.772%
9	20.562	1683	1690	1707	rBV	3710848	8589317	100.00%	18.116%
10	21.067	1738	1745	1765	rBV2	45767	233516	2.72%	0.493%
11	22.793	1926	1933	1941	rBV	232503	499871	5.82%	1.054%
12	24.987	2163	2172	2210	rBV2	2990330	8501933	98.98%	17.931%
13	26.888	2374	2379	2386	rBV	95299	204536	2.38%	0.431%
14	28.659	2566	2572	2579	rBV	62344	139827	1.63%	0.295%
15	30.064	2719	2725	2738	rBV	413437	887965	10.34%	1.873%
16	30.248	2738	2745	2752	rVB	51714	124064	1.44%	0.262%
17	31.716	2900	2905	2912	rVB	38852	94416	1.10%	0.199%
18	33.020	3040	3047	3067	rBV2	2145664	5829561	67.87%	12.295%
19	35.572	3320	3325	3333	rBV	29441	93258	1.09%	0.197%
20	37.114	3485	3493	3521	rBV2	1196511	4031049	46.93%	8.502%

Sum of corrected areas: 47413952

30608.D GCMS0103.M Thu Jan 10 15:07:34 2002

LSC Report - Integrated Chromatogram

File : M:\INSTRU~1\209\DATA\DEC2701\30608.D
Operator : 209 GALOS
Acquired : 28 Dec 2001 8:16 pm using AcqMethod GCMS1126
Instrument : GC/MS 209
Sample Name: GC/MS SCAN 12/17
Misc Info :
Vial Number: 32
Quant File :GCMS1126.RES (RTE Integrator)



30608.D GCMS0103.M Thu Jan 10 15:07:41 2002

Library Search Compound Report

Data File : M:\INSTRU~1\209\DATA\DEC2701\30608.D
Acq On : 28 Dec 2001 8:16 pm
Sample : GC/MS SCAN 12/17
Misc :
MS Integration Params: LSCINT.P

Vial: 32
Operator: 209 GALOS
Inst : GC/MS 209
Multiplr: 1.00

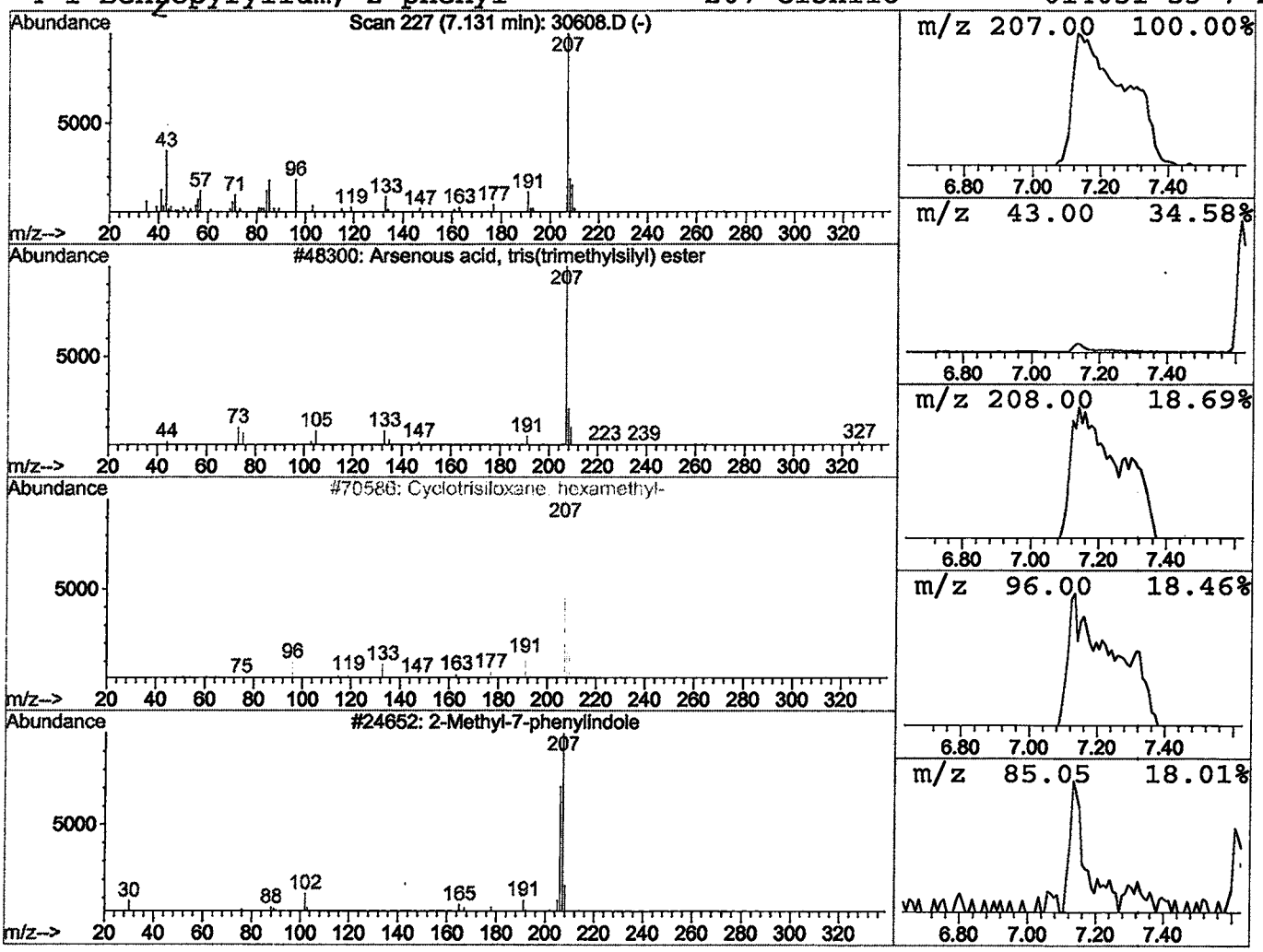
Quant Method : C:\HPCHEM\1\METHODS\GCMS1126.M (RTE Integrator)
Title : 209 625/TCLP calibration table
Library : C:\DATABASE\NBS75K.L

****

Peak Number 1 Arsenous acid, tris(trimethyls Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.13	0.41 ug/l	130123	1,4-Dichlorobenzene-d4	11.68

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Arsenous acid, tris(trimethylsilyl)	342	C9H27AsO3Si3	055429-29-3	42
2	Cyclotrisiloxane, hexamethyl-	222	C6H18O3Si3	000541-05-9	38
3	2-Methyl-7-phenylindole	207	C15H13N	000000-00-0	33
4	1-Benzopyrylium, 2-phenyl-	207	C15H11O	014051-53-7	28



Library Search Compound Report

Data File : M:\INSTRU~1\209\DATA\DEC2701\30608.D
Acq On : 28 Dec 2001 8:16 pm
Sample : GC/MS SCAN 12/17
Misc :
MS Integration Params: LSCINT.P

Vial: 32
Operator: 209 GALOS
Inst : GC/MS 209
Multiplr: 1.00

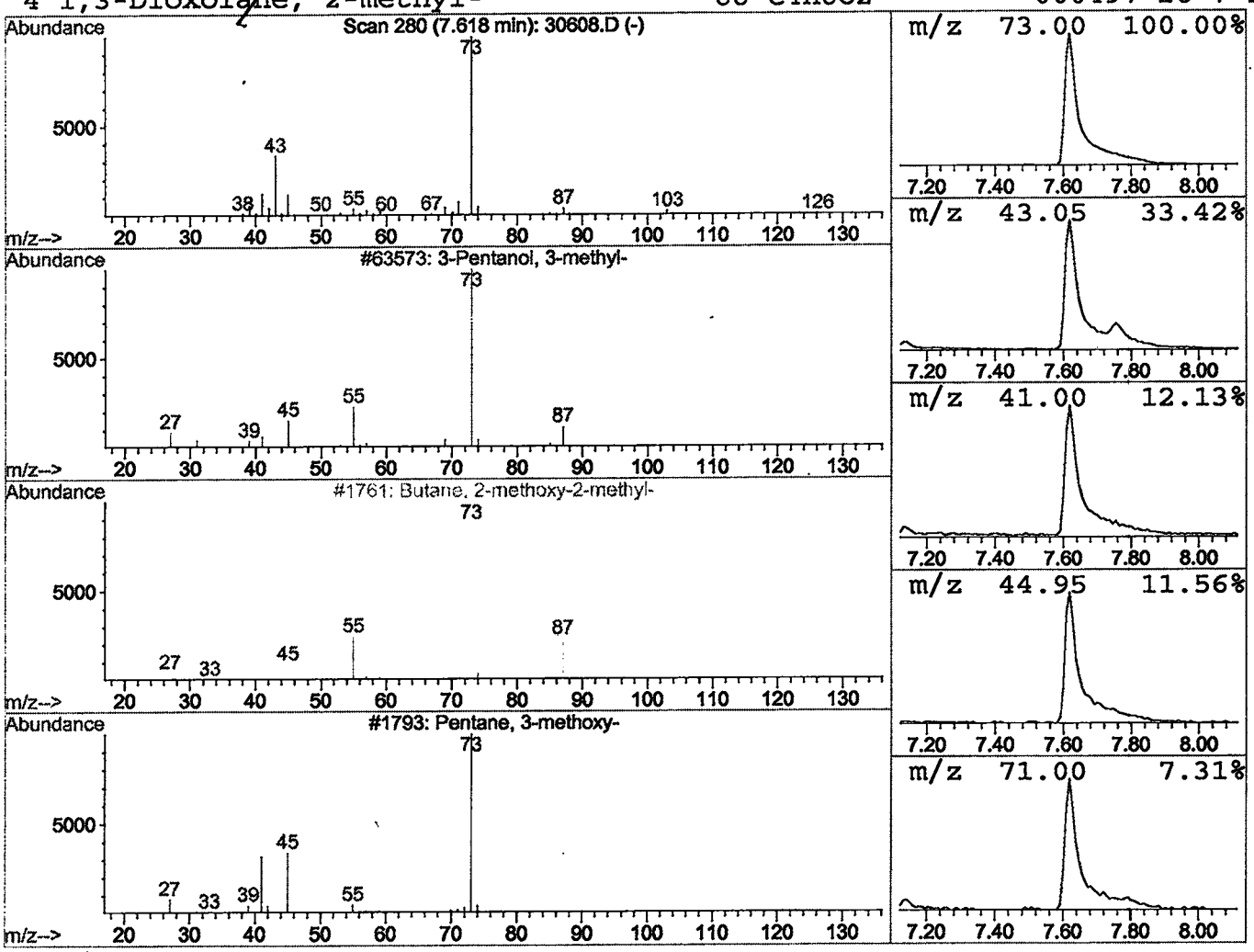
Quant Method : C:\HPCHEM\1\METHODS\GCMS1126.M (RTE Integrator)
Title : 209 625/TCLP calibration table
Library : C:\DATABASE\NBS75K.L

****

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.62	3.52 ug/l	1116870	1,4-Dichlorobenzene-d4	11.68

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	3-Pentanol, 3-methyl-	102	C6H14O	000077-74-7	28
2	Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	23
3	Pentane, 3-methoxy-	102	C6H14O	036839-67-5	17
4	1,3-Dioxolane, 2-methyl-	88	C4H8O2	000497-26-7	9



Library Search Compound Report

Data File : M:\INSTRU~1\209\DATA\DEC2701\30608.D Vial: 32
Acq On : 28 Dec 2001 8:16 pm Operator: 209 GALOS
Sample : GC/MS SCAN 12/17 Inst : GC/MS 209
Misc : Multiplr: 1.00
MS Integration Params: LSCINT.P

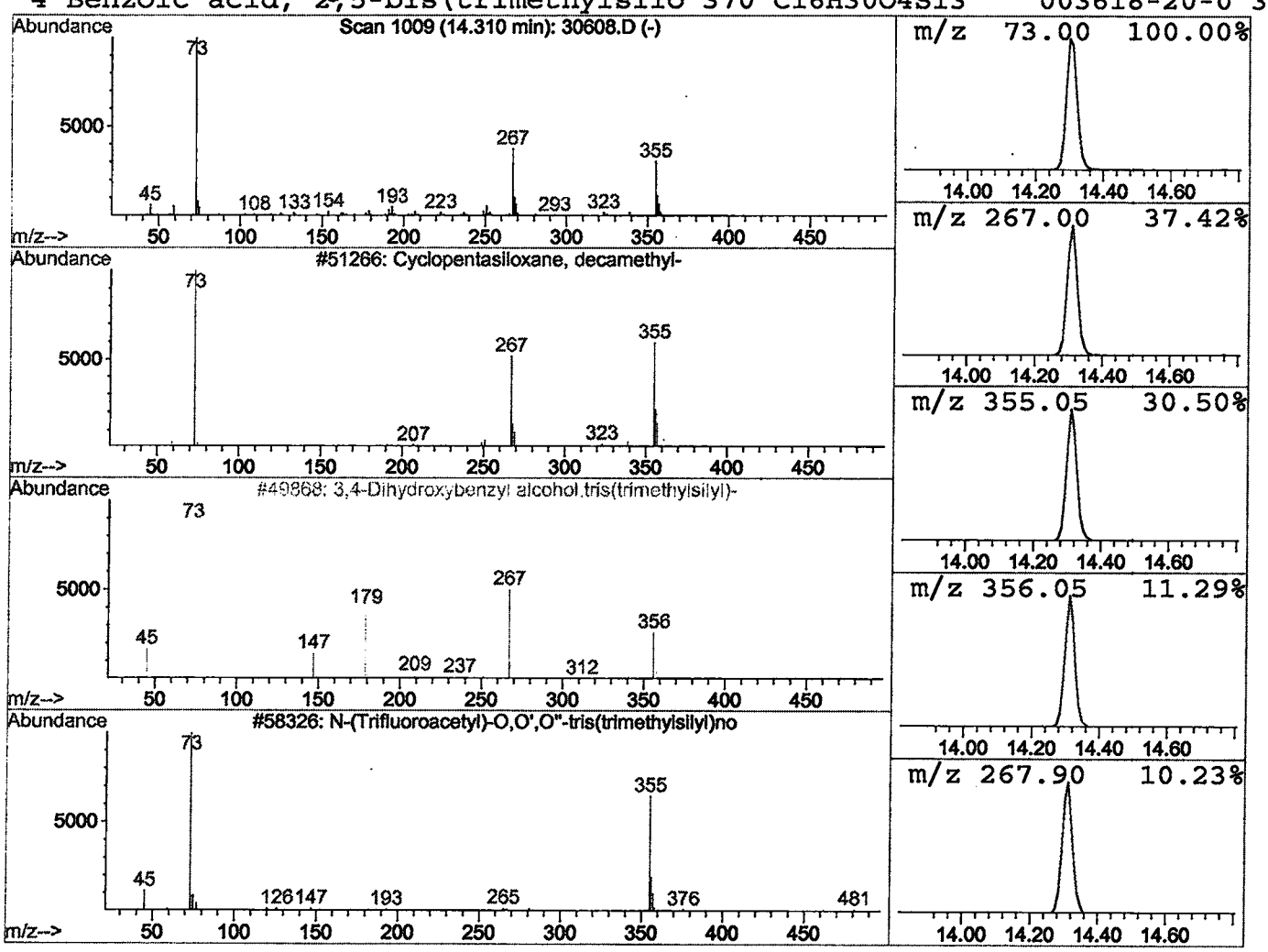
Quant Method : C:\HPCHEM\1\METHODS\GCMS1126.M (RTE Integrator)
Title : 209 625/TCLP calibration table
Library : C:\DATABASE\NBS75K.L

****

Peak Number 3 Cyclopentasiloxane, decamethyl Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.31	1.31 ug/l	518333	Naphthalene-d8	15.28

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclopentasiloxane, decamethyl-	370	C10H30O5Si5	000541-02-6	72
2	3,4-Dihydroxybenzyl alcohol,tris(tr	356	C16H32O3Si3	068595-79-9	38
3	N-(Trifluoroacetyl)-O,O',O''-tris(t	481	C19H34F3NO4Si3	000000-00-0	37
4	Benzoic acid, 2,5-bis(trimethylsilo	370	C16H30O4Si3	003618-20-0	37



Library Search Compound Report

Data File : M:\INSTRU~1\209\DATA\DEC2701\30608.D Vial: 32
Acq On : 28 Dec 2001 8:16 pm Operator: 209 GALOS
Sample : GC/MS SCAN 12/17 Inst : GC/MS 209
Misc : Multiplr: 1.00
MS Integration Params: LSCINT.P

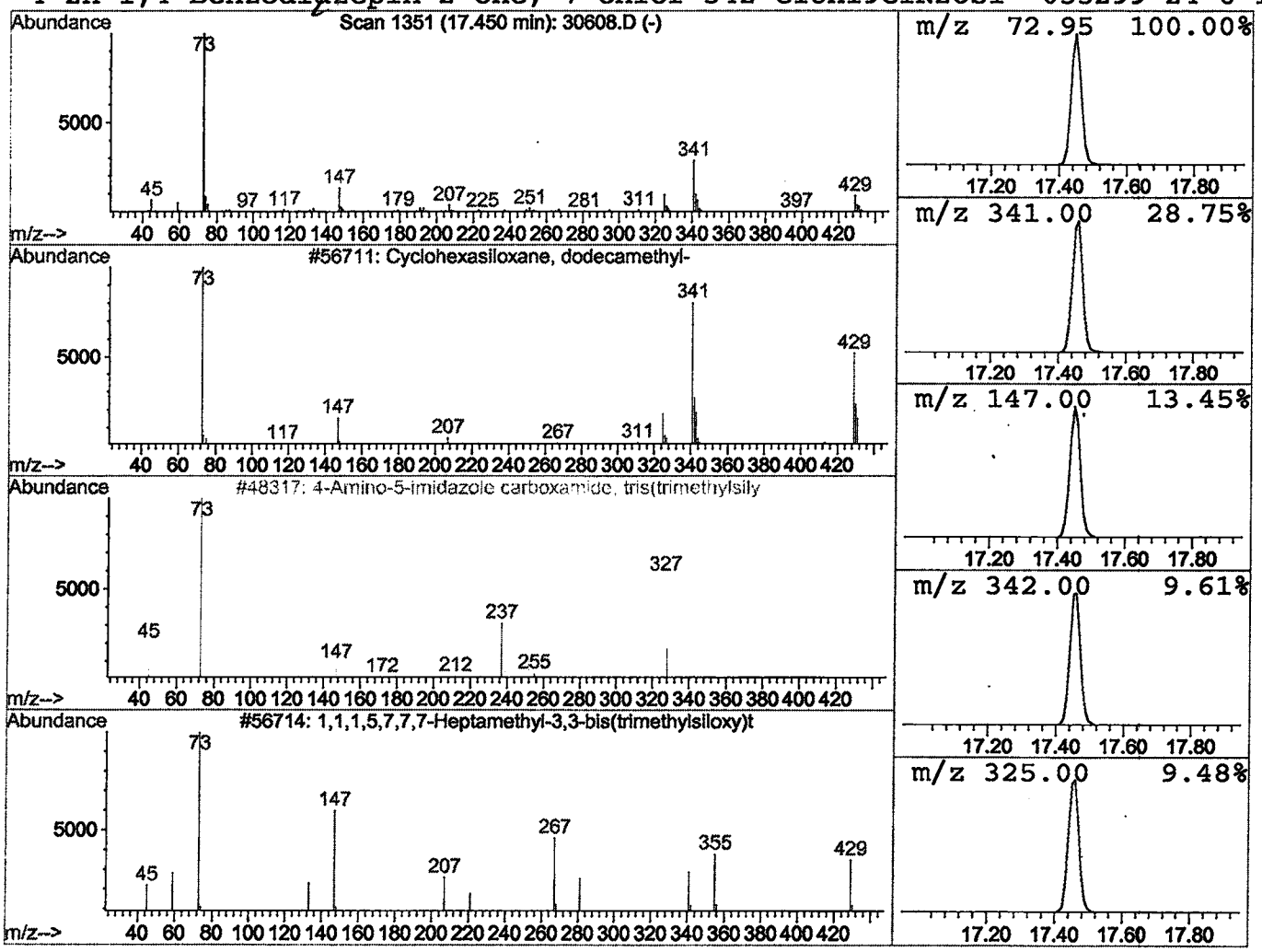
Quant Method : C:\HPCHEM\1\METHODS\GCMS1126.M (RTE Integrator)
Title : 209 625/TCLP calibration table
Library : C:\DATABASE\NBS75K.L

****

Peak Number 4 Cyclohexasiloxane, dodecamethy Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.45	3.01 ug/l	1193220	Naphthalene-d8	15.28

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclohexasiloxane, dodecamethyl-	444	C12H36O6Si6	000540-97-6	38
2	4-Amino-5-imidazole carboxamide, tr	342	C13H30N4OSi3	000000-00-0	32
3	1,1,1,5,7,7,7-Heptamethyl-3,3-bis(t	444	C13H40O5Si6	038147-00-1	25
4	2H-1,4-Benzodiazepin-2-one, 7-chlor	342	C18H19ClN2OSi	055299-24-6	12



Library Search Compound Report

Data File : M:\INSTRU~1\209\DATA\DEC2701\30608.D Vial: 32
Acq On : 28 Dec 2001 8:16 pm Operator: 209 GALOS
Sample : GC/MS SCAN 12/17 Inst : GC/MS 209
Misc : Multiplr: 1.00
MS Integration Params: LSCINT.P

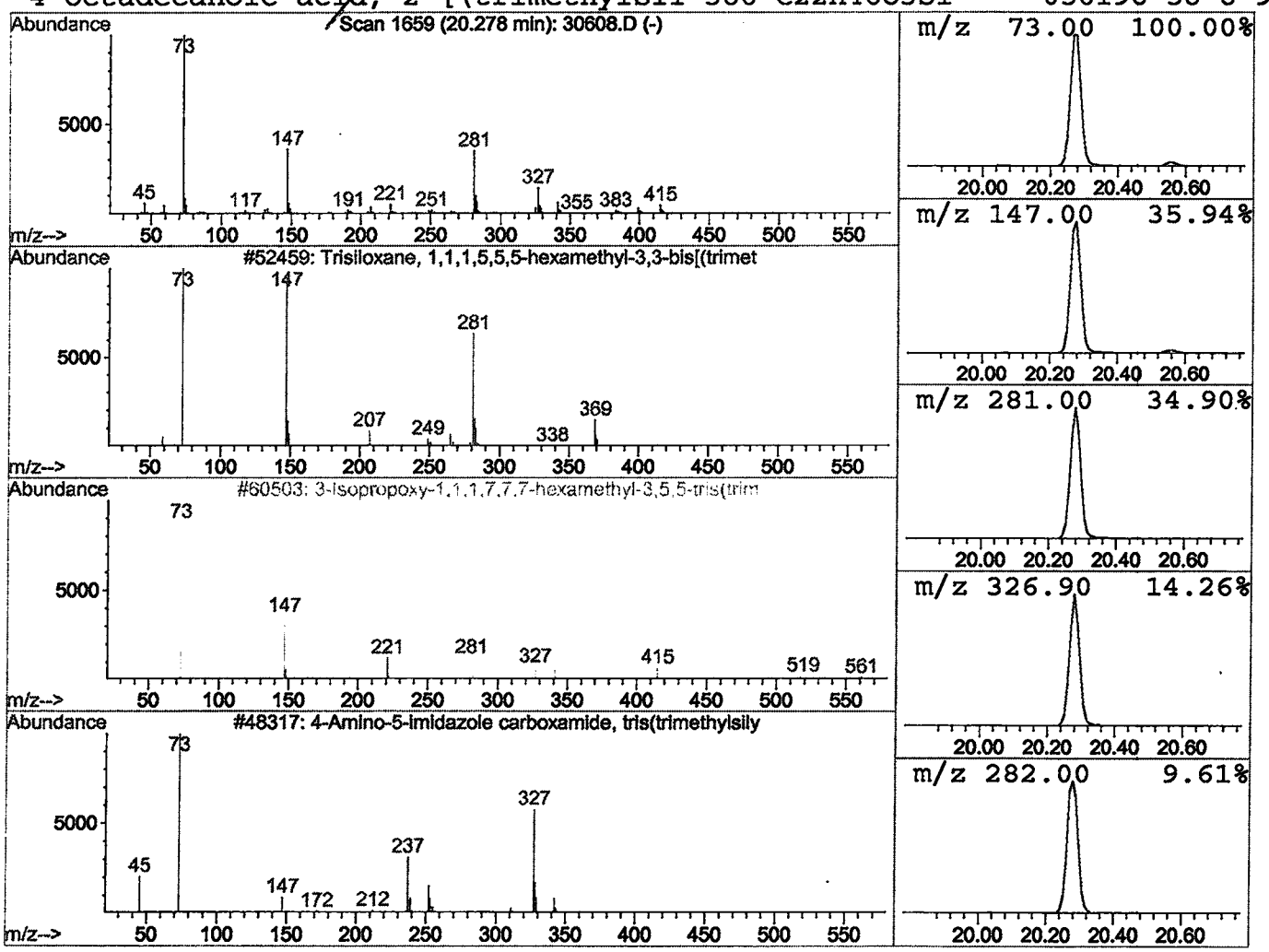
Quant Method : C:\HPCHEM\1\METHODS\GCMS1126.M (RTE Integrator)
Title : 209 625/TCLP calibration table
Library : C:\DATABASE\NBS75K.L

****

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	1.96 ug/l	839991	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	4-Amino-5-imidazole carboxamide, tr	342	C13H30N4OSi3	000000-00-0	10
4	Octadecanoic acid, 2-[(trimethylsil	386	C22H46O3Si	056196-58-8	9



Library Search Compound Report

Data File : M:\INSTRU~1\209\DATA\DEC2701\30608.D
Acq On : 28 Dec 2001 8:16 pm
Sample : GC/MS SCAN 12/17
Misc :
MS Integration Params: LSCINT.P

Vial: 32
Operator: 209 GALOS
Inst : GC/MS 209
Multiplr: 1.00

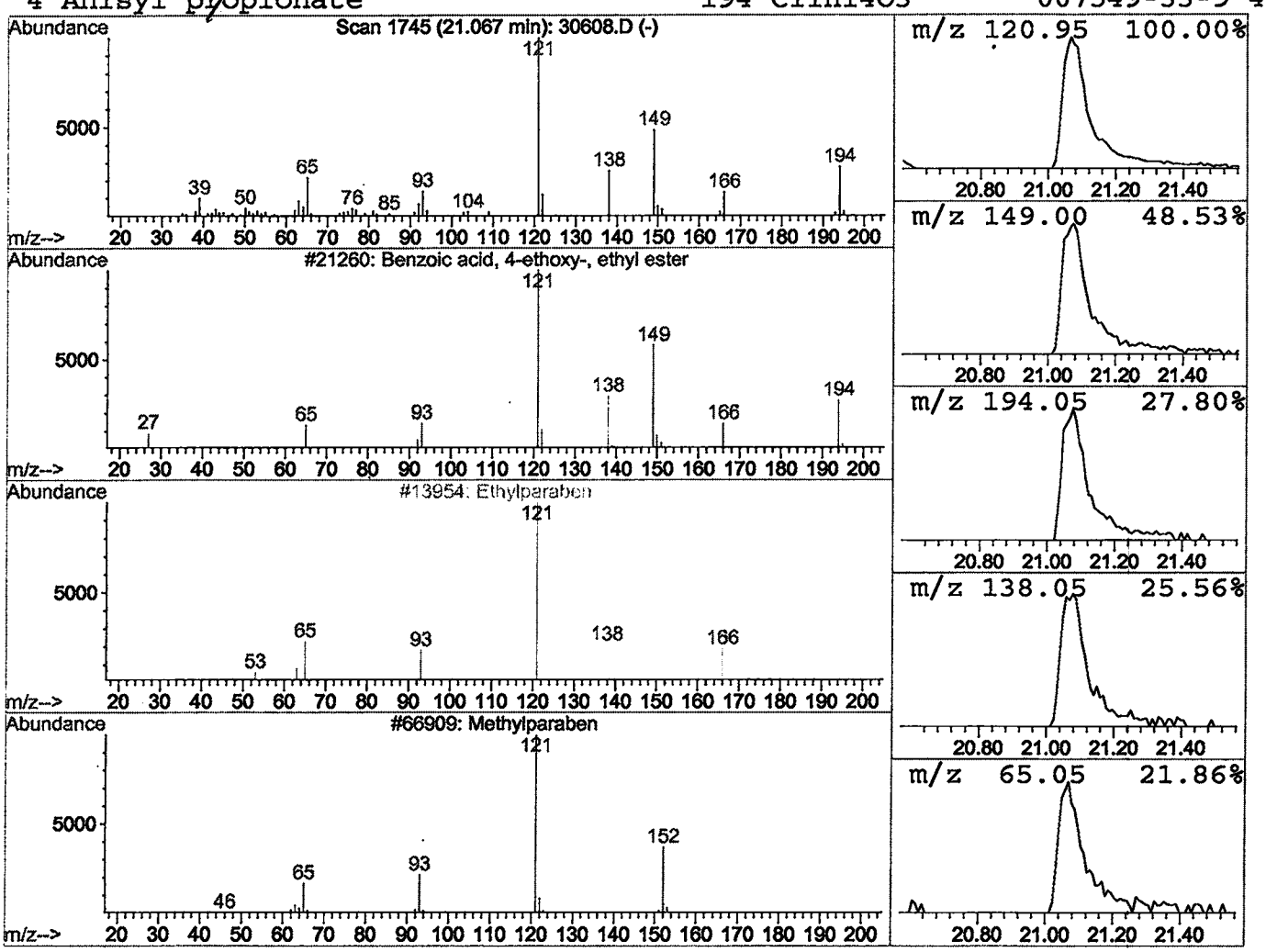
Quant Method : C:\HPCHEM\1\METHODS\GCMS1126.M (RTE Integrator)
Title : 209 625/TCLP calibration table
Library : C:\DATABASE\NBS75K.L

****

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.07	0.54 ug/l	233516	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	94
2	Ethylparaben	166	C9H10O3	000120-47-8	53
3	Methylparaben	152	C8H8O3	000099-76-3	43
4	Anisyl propionate	194	C11H14O3	007549-33-9	43



Library Search Compound Report

Data File : M:\INSTRU~1\209\DATA\DEC2701\30608.D Vial: 32
Acq On : 28 Dec 2001 8:16 pm Operator: 209 GALOS
Sample : GC/MS SCAN 12/17 Inst : GC/MS 209
Misc : Multiplr: 1.00
MS Integration Params: LSCINT.P

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

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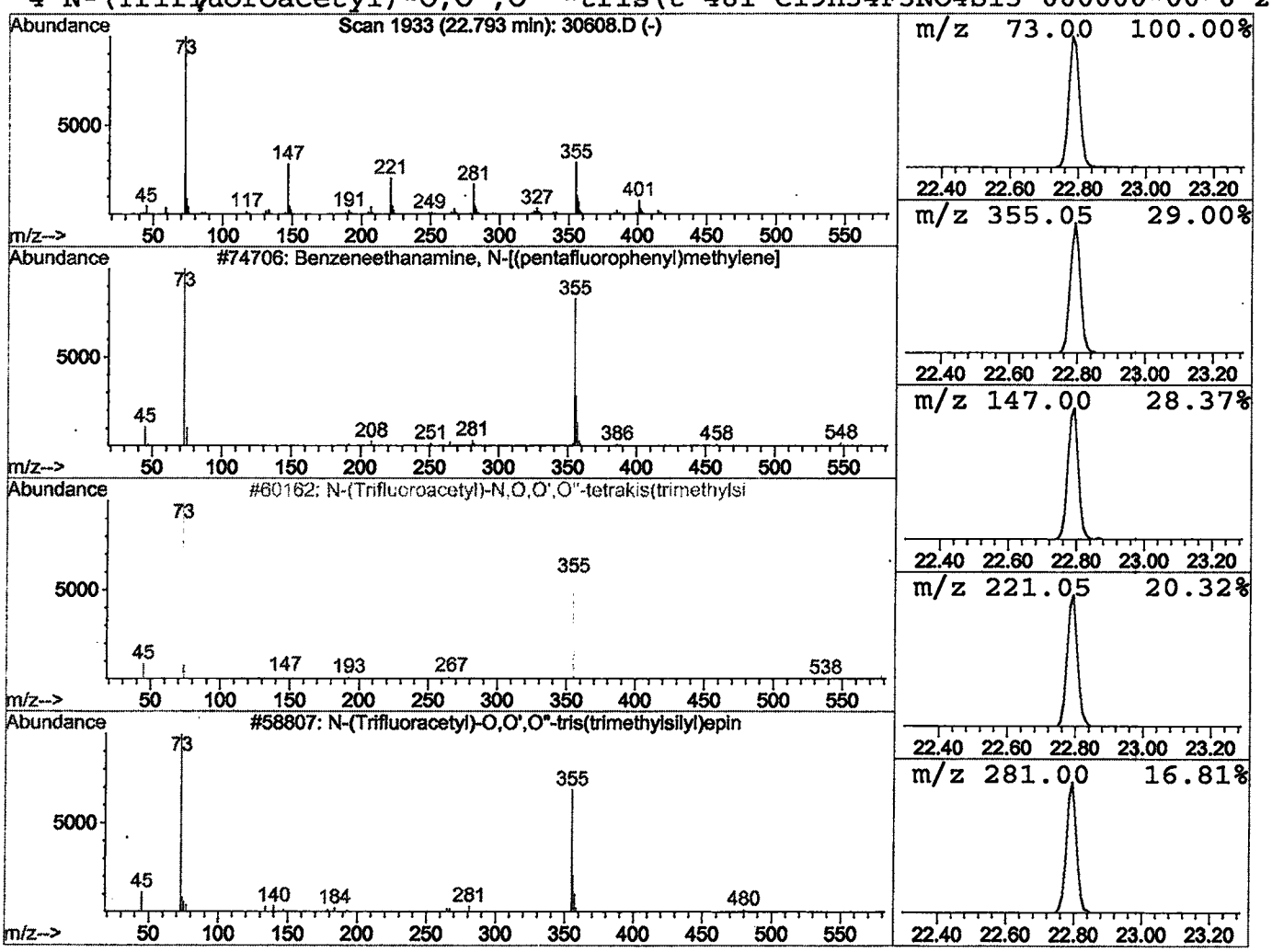
Peak Number 7 Benzeneethanamine, N-[(pentafl Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.79	1.18 ug/l	499871	Phenanthrene-d10	24.99

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzeneethanamine, N-[(pentafluorop	563	C24H34F5NO3Si3	055429-13-5	25
2	N-(Trifluoroacetyl)-N,O,O',O''-tetr	553	C22H42F3NO4Si4	000000-00-0	22
3	N-(Trifluoracetyl)-O,O',O"-tris(tri	495	C20H36F3NO4Si3	000000-00-0	22
4	N-(Trifluoroacetyl)-O,O',O''-tris(t	481	C19H34F3NO4Si3	000000-00-0	22



Library Search Compound Report

Data File : M:\INSTRU~1\209\DATA\DEC2701\30608.D
Acq On : 28 Dec 2001 8:16 pm
Sample : GC/MS SCAN 12/17
Misc :
MS Integration Params: LSCINT.P

Vial: 32
Operator: 209 GALOS
Inst : GC/MS 209
Multiplr: 1.00

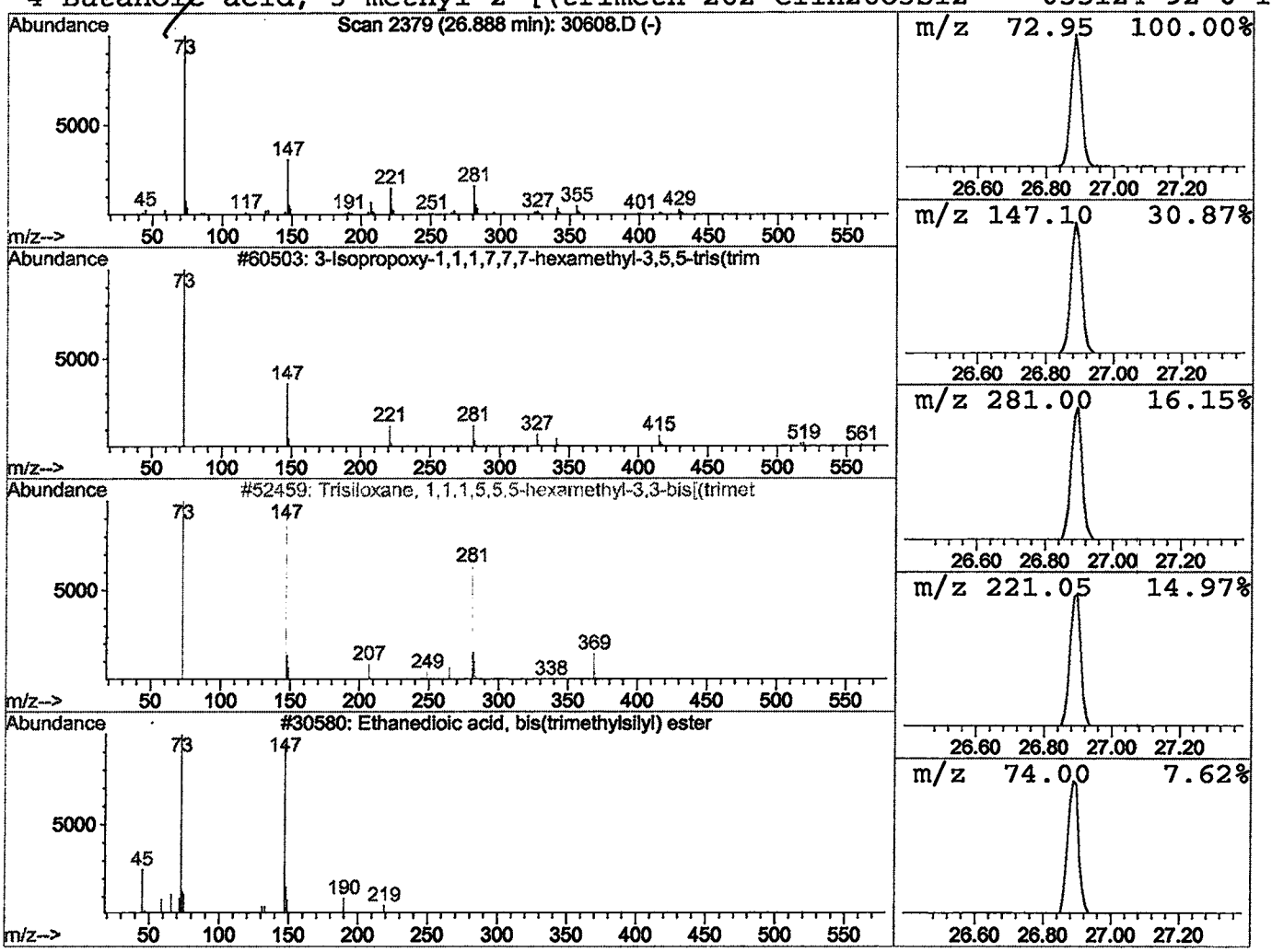
Quant Method : C:\HPCHEM\1\METHODS\GCMS1126.M (RTE Integrator)
Title : 209 625/TCLP calibration table
Library : C:\DATABASE\NBS75K.L

****

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
26.89	0.48 ug/l	204536	Phenanthrene-d10	24.99

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	3-Isopropoxy-1,1,1,7,7,7-hexamethyl	576	C18H52O7Si7	071579-69-6	56
2	Trisiloxane, 1,1,1,5,5,5-hexamethyl	384	C12H36O4Si5	003555-47-3	40
3	Ethanedioic acid, bis(trimethylsily	234	C8H18O4Si2	018294-04-7	25
4	Butanoic acid, 3-methyl-2-[(trimeth	262	C11H26O3Si2	055124-92-0	17



Library Search Compound Report

Data File : M:\INSTRU~1\209\DATA\DEC2701\30608.D Vial: 32
Acq On : 28 Dec 2001 8:16 pm Operator: 209 GALOS
Sample : GC/MS SCAN 12/17 Inst : GC/MS 209
Misc : Multiplr: 1.00
MS Integration Params: LSCINT.P

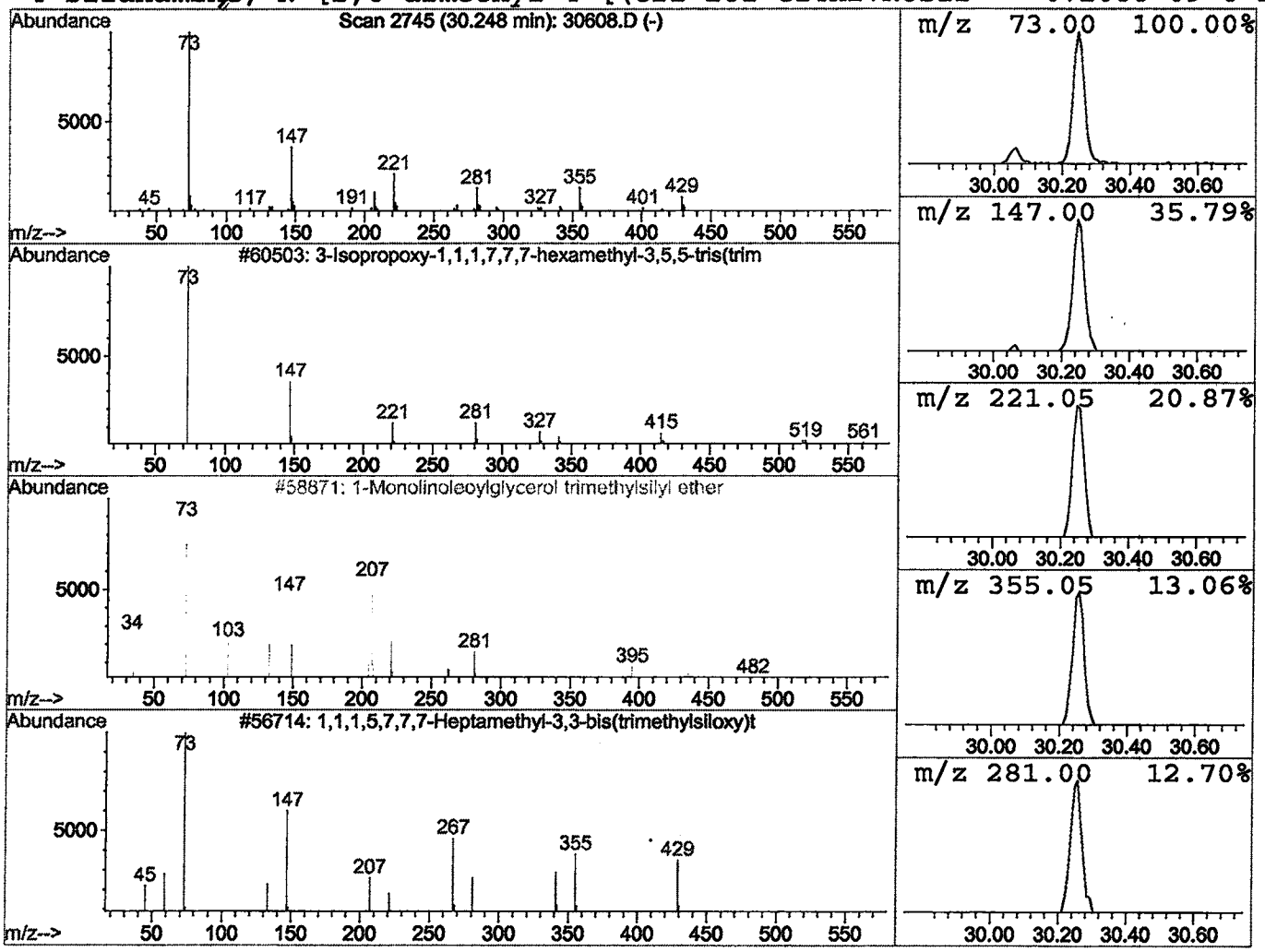
Quant Method : C:\HPCHEM\1\METHODS\GCMS1126.M (RTE Integrator)
Title : 209 625/TCLP calibration table
Library : C:\DATABASE\NBS75K.L

****

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
30.25	0.43 ug/l	124064	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	28
2	1-Monolinoleoylglycerol trimethylsi	498	C27H54O4Si2	054284-45-6	28
3	1,1,1,5,7,7,7-Heptamethyl-3,3-bis(t	444	C13H40O5Si6	038147-00-1	23
4	Silanamine, N-[2,6-dimethyl-4-[(tri	281	C14H27NOSi2	072088-09-6	22



Library Search Compound Report

Data File : M:\INSTRU~1\209\DATA\DEC2701\30608.D
Acq On : 28 Dec 2001 8:16 pm
Sample : GC/MS SCAN 12/17
Misc :
MS Integration Params: LSCINT.P

Vial: 32
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

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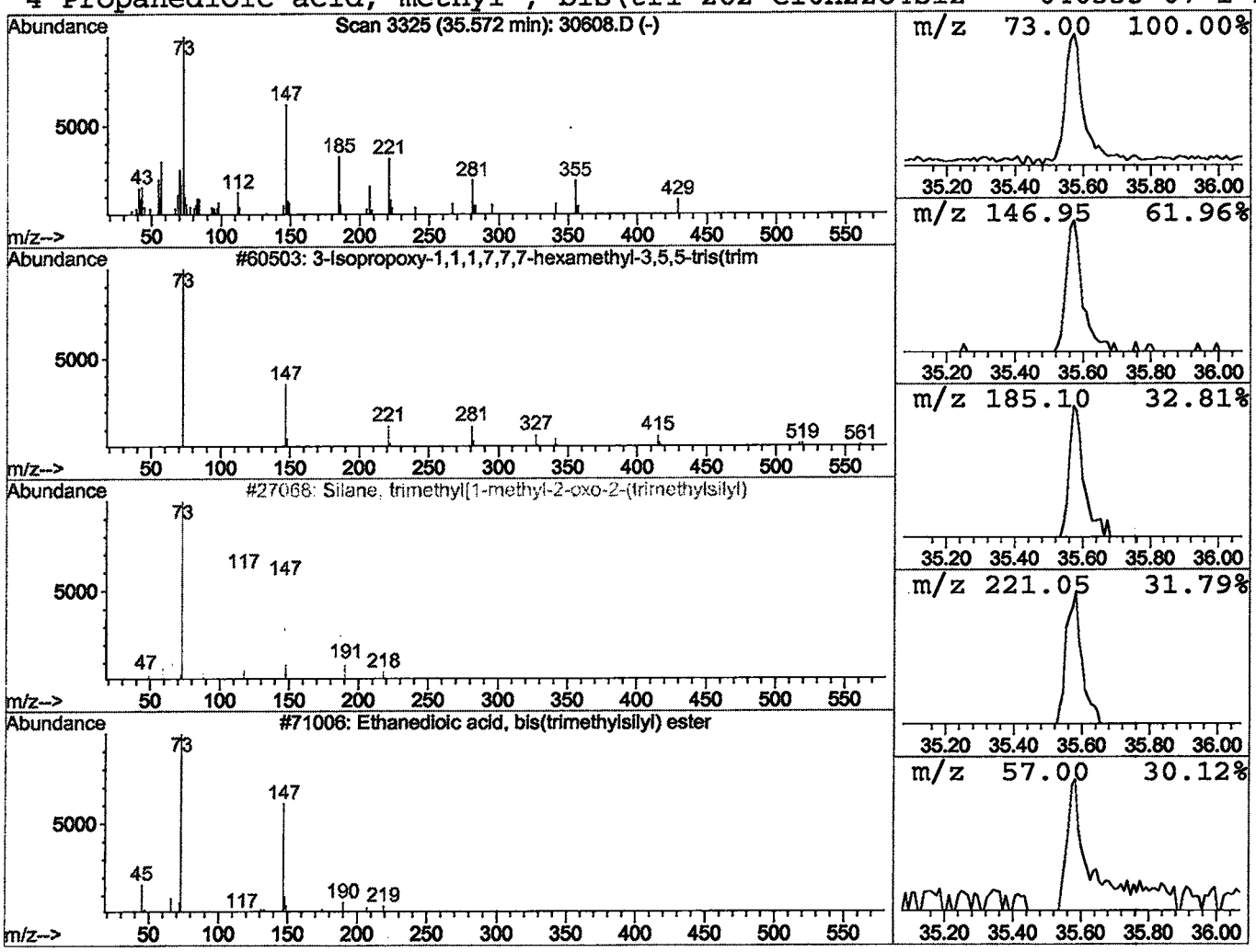
Peak Number 10 3-Isopropoxy-1,1,1,7,7,7-hexam Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
35.57	0.46 ug/l	93258	Perylene-d12	37.11

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	3-Isopropoxy-1,1,1,7,7,7-hexamethyl	576	C18H52O7Si7	071579-69-6	17
2	Silane, trimethyl[1-methyl-2-oxo-2-	218	C9H22O2Si2	055255-93-1	10
3	Ethanedioic acid, bis(trimethylsily	234	C8H18O4Si2	018294-04-7	10
4	Propanedioic acid, methyl-, bis(tri	262	C10H22O4Si2	040333-07-1	10



Tentatively Identified Compound (LSC) summary

Operator ID: 209 GALOS Date Acquired: 28 Dec 2001 8:16 pm
Data File: M:\INSTRU~1\209\DATA\DEC2701\30608.D
Name: GC/MS SCAN 12/17
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
Arsenous acid, tris(	7.13	0.4	ug/l	130123	ISTD01	11.68	6351980	20.0
3-Pentanol, 3-methyl	7.62	3.5	ug/l	1116870	ISTD01	11.68	6351980	20.0
Cyclopentasiloxane,	14.31	1.3	ug/l	518333	ISTD02	15.28	7936870	20.0
Cyclohexasiloxane, d	17.45	3.0	ug/l	1193220	ISTD02	15.28	7936870	20.0
Trisiloxane, 1,1,1,5	20.28	2.0	ug/l	839991	ISTD03	20.56	8589320	20.0
Benzoic acid, 4-etho	21.07	0.5	ug/l	233516	ISTD03	20.56	8589320	20.0
Benzeneethanamine, N	22.79	1.2	ug/l	499871	ISTD04	24.99	8501930	20.0
3-Isopropoxy-1,1,1,7	26.89	0.5	ug/l	204536	ISTD04	24.99	8501930	20.0
3-Isopropoxy-1,1,1,7	30.25	0.4	ug/l	124064	ISTD05	33.02	5829560	20.0
3-Isopropoxy-1,1,1,7	35.57	0.5	ug/l	93258	ISTD06	37.11	4031050	20.0

30608.D GCMS0103.M Thu Jan 10 15:08:37 2002