User: Vinci, David L
Westchester Dept. of Labs & Research
Date: 01/25/2002 Time: 12:20:13

Sample: AD30713
Analysis: $GCMS GCMS Scan For Unknowns
Units: ug
Started: 1/25/02 12:19 Ended: 1/25/02 12:19 Analyst: DLV
Page: 1

	Component Name	Viol	Result	Qualifier	MDL
1	Other unknown compounds		.		
2	Surr_2,4-DDD		99		5.0

Comments for Sample AD30713 - Analysis $GCMS

Westchester Dept. of Labs & Research
User: Vinci, David L
Date: 01-25-2002 Time: 12:20:27

The GCMS scan, from 35 to 550 amu, does not reveal the presence of unusual compounds.
The recoveries for 2 of the 43 compounds in the LCS Standard were above their acceptance ranges.

Quantitation Report (QT Reviewed)

Data File : C:\HPCHEM\1\DATA\DEC3101\30713.D Vial: 28
Acq On : 1 Jan 2002 3:07 pm Operator: 209 GALOS
Sample : gcms scan 12/19 Inst : GC/MS 209
Misc : Multiplr: 1.00
MS Integration Params: GOOFY.P
Quant Time: Jan 18 11:00 2002 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	923092	20.00	ug/l	-0.02
10) Naphthalene-d8	15.28	136	3506560	20.00	ug/l	-0.02
17) Acenaphthene-d10	20.56	164	1808793	20.00	ug/l	-0.01
29) Phenanthrene-d10	24.99	188	2623578m	20.00	ug/l	0.00
38) Chrysene-d12	33.02	240	1571973	20.00	ug/l	-0.01
44) Perylene-d12	37.11	264	841459	20.00	ug/l	0.00
System Monitoring Compounds						
37) 2,4-DDD	30.06	235	142523	4.94	ug/mL	0.00
Spiked Amount 5.000			Recovery = 98.80%			
Target Compounds						Qvalue
2) N-nitroso-dimethyl amine	5.38	74	620	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	15.34	128	606	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	251	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	21.39	165	209	N.D.		
25) Diethylphthalate	22.12	149	2639	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)

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Acq On : 1 Jan 2002 3:07 pm Operator: 209 GALOS
Sample : gcms scan 12/19 Inst : GC/MS 209
Misc : Multiplr: 1.00
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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	24.98	178	856	N.D.	
34)	Anthracene	24.98	178	856	N.D.	
35)	Di-n-butylphthalate	27.03	149	8386	N.D.	
36)	Fluoranthene	0.00	202	0	N.D.	
39)	Pyrene	0.00	202	0	N.D.	
40)	Butylbenzylphthalate	31.51	149	424	N.D.	
41)	Benzo(a)anthracene	33.03	228	5698	N.D.	
42)	Bis(2-ethylhexyl)phthalate	33.30	149	4659	N.D.	
43)	Chrysene	33.09	228	2579	N.D.	
45)	Di-n-Octylphthalate	35.04	149	747	N.D.	
46)	Benzo(b)fluoranthene	36.11	252	746	N.D.	
47)	Benzo(k)fluoranthene	36.17	252	1802	N.D.	
48)	Benzo(a)pyrene	37.11	252	4017	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

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Sample : gcms scan 12/19
Misc :
MS Integration Params: GOOFY.P
Quant Time: Jan 18 11:00 2002

Vial: 28
Operator: 209 GALOS
Inst : GC/MS 209
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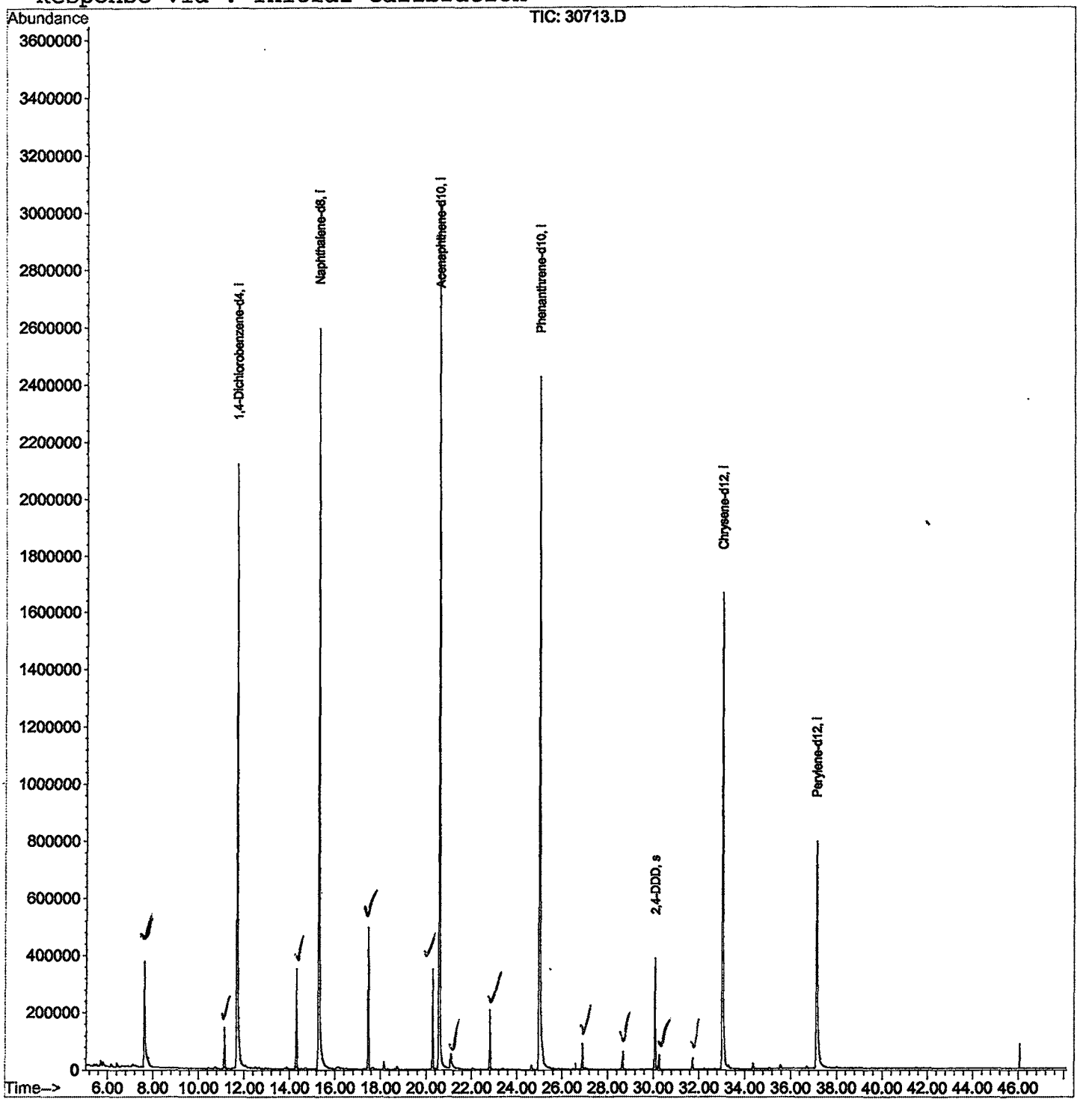
Quant Results File: GCMS1126.RES

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\DEC3101\30713.D Vial: 28
Acq On : 1 Jan 2002 3:07 pm Operator: 209 GALOS
Sample : gcms scan 12/19 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

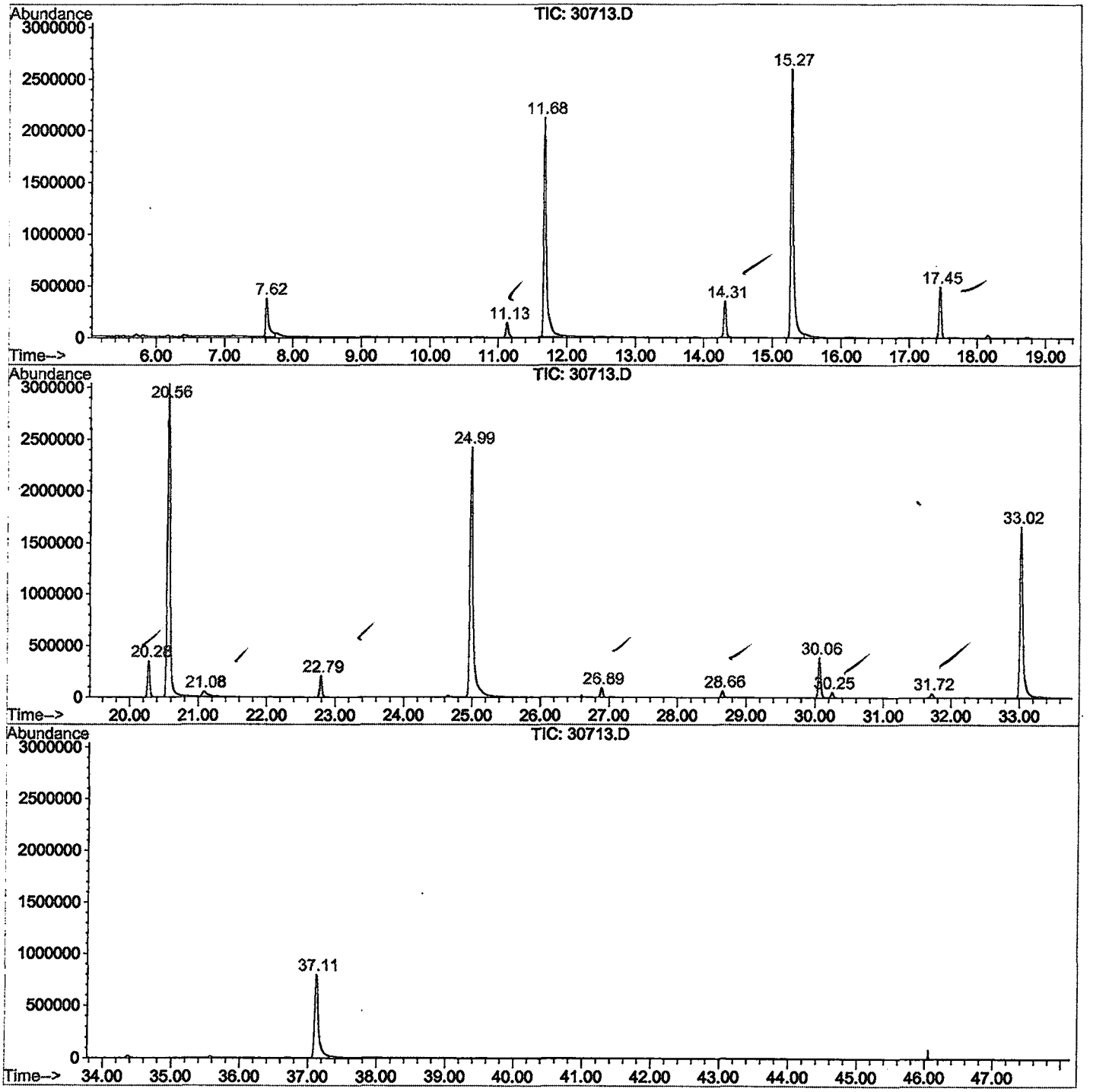
peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	peak area	peak % max.	% of total
1	7.618	275	280	294	rBV	370092	1118122	14.13%	2.591%
2	11.134	657	663	676	rVB	143389	363659	4.60%	0.843%
3	11.676	717	722	751	rBV	2115596	5713713	72.21%	13.243%
4	14.310	999	1009	1020	rBV	349192	816347	10.32%	1.892%
5	15.274	1109	1114	1134	rBV	2594009	7027208	88.81%	16.287%
6	17.450	1344	1351	1362	rBV	493724	1161248	14.68%	2.691%
7	20.277	1651	1659	1666	rBV	350524	762328	9.63%	1.767%
8	20.562	1683	1690	1711	rBV	3036379	7912596	100.00%	18.339%
9	21.076	1740	1746	1762	rBV2	49362	247350	3.13%	0.573%
10	22.793	1925	1933	1944	rBV	206736	464199	5.87%	1.076%
11	24.987	2164	2172	2217	rBV2	2425289	7840912	99.09%	18.173%
12	26.887	2373	2379	2386	rBV	88748	201540	2.55%	0.467%
13	28.659	2565	2572	2579	rBV	62156	142441	1.80%	0.330%
14	30.064	2719	2725	2734	rBV	387182	897410	11.34%	2.080%
15	30.247	2740	2745	2755	rVB	51541	123012	1.55%	0.285%
16	31.716	2898	2905	2912	rBV	39741	102071	1.29%	0.237%
17	33.020	3040	3047	3074	rBV2	1663525	5195459	65.66%	12.041%
18	37.114	3486	3493	3515	rBV2	792152	3057006	38.63%	7.085%

Sum of corrected areas: 43146621

30713.D GCMS1126.M Wed Jan 23 08:46:01 2002

LSC Report - Integrated Chromatogram

File : C:\HPCHEM\1\DATA\DEC3101\30713.D
Operator : 209 GALOS
Acquired : 1 Jan 2002 3:07 pm using AcqMethod GCMS1126
Instrument : GC/MS 209
Sample Name: gcms scan 12/19
Misc Info :
Vial Number: 28
Quant File :GCMS1126.RES (RTE Integrator)



30713.D GCMS1126.M Wed Jan 23 08:46:07 2002

Library Search Compound Report

Data File : C:\HPCHEM\1\DATA\DEC3101\30713.D
Acq On : 1 Jan 2002 3:07 pm
Sample : gcms scan 12/19
Misc :
MS Integration Params: LSCINT.P

Vial: 28
Operator: 209 GALOS
Inst : GC/MS 209
Multiplr: 1.00

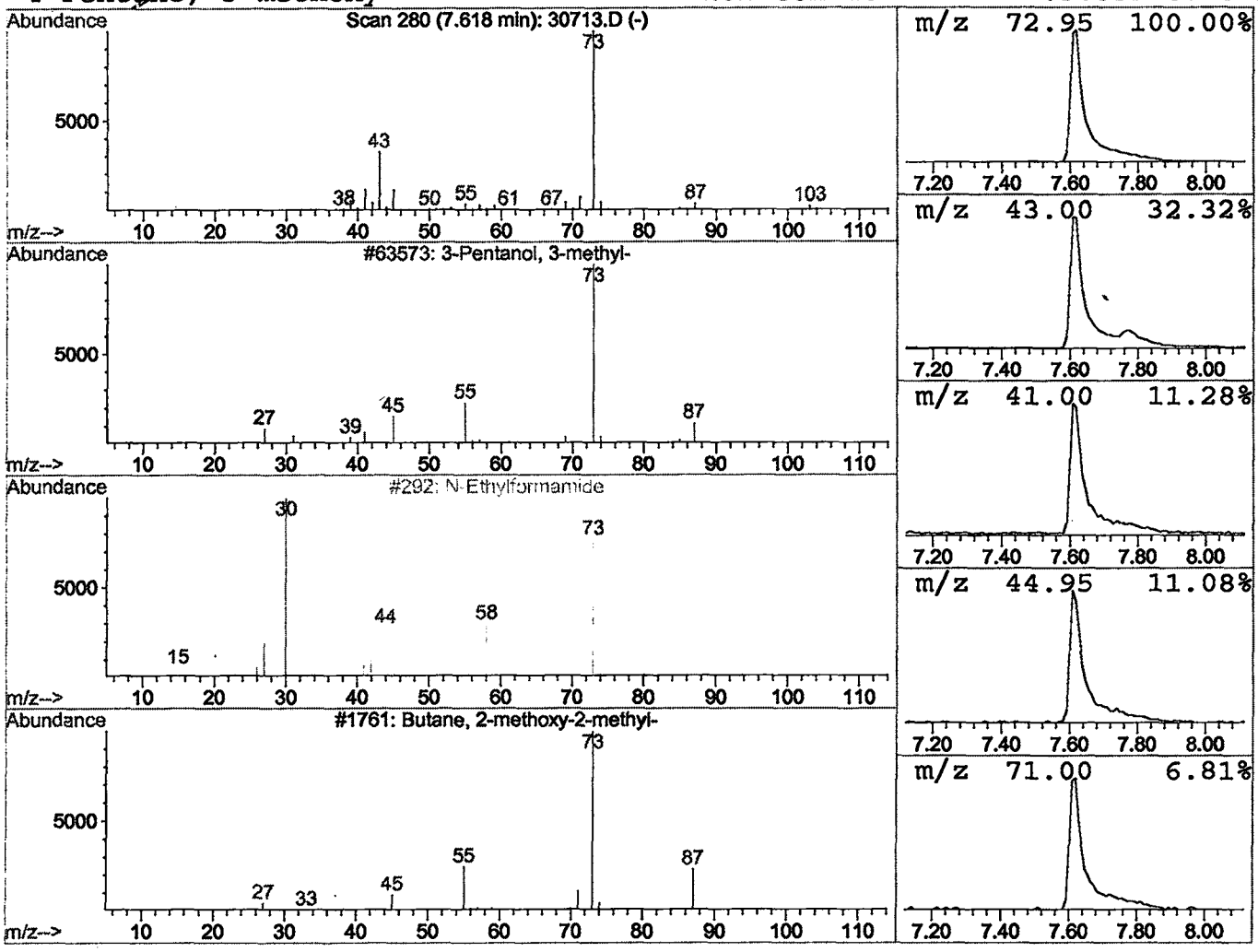
Quant Method : C:\HPCHEM\1\METHODS\GCMS1126.M (RTE Integrator)
Title : 209 625/TCLP calibration table
Library : C:\DATABASE\NBS75K.L

****

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.62	3.91 ug/l	1118120	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	33
2	N-Ethylformamide	73	C3H7NO	000627-45-2	9
3	Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	9
4	Pentane, 3-methoxy-	102	C6H14O	036839-67-5	9



Library Search Compound Report

Data File : C:\HPCHEM\1\DATA\DEC3101\30713.D
Acq On : 1 Jan 2002 3:07 pm
Sample : gcms scan 12/19
Misc :
MS Integration Params: LSCINT.P

Vial: 28
Operator: 209 GALOS
Inst : GC/MS 209
Multiplr: 1.00

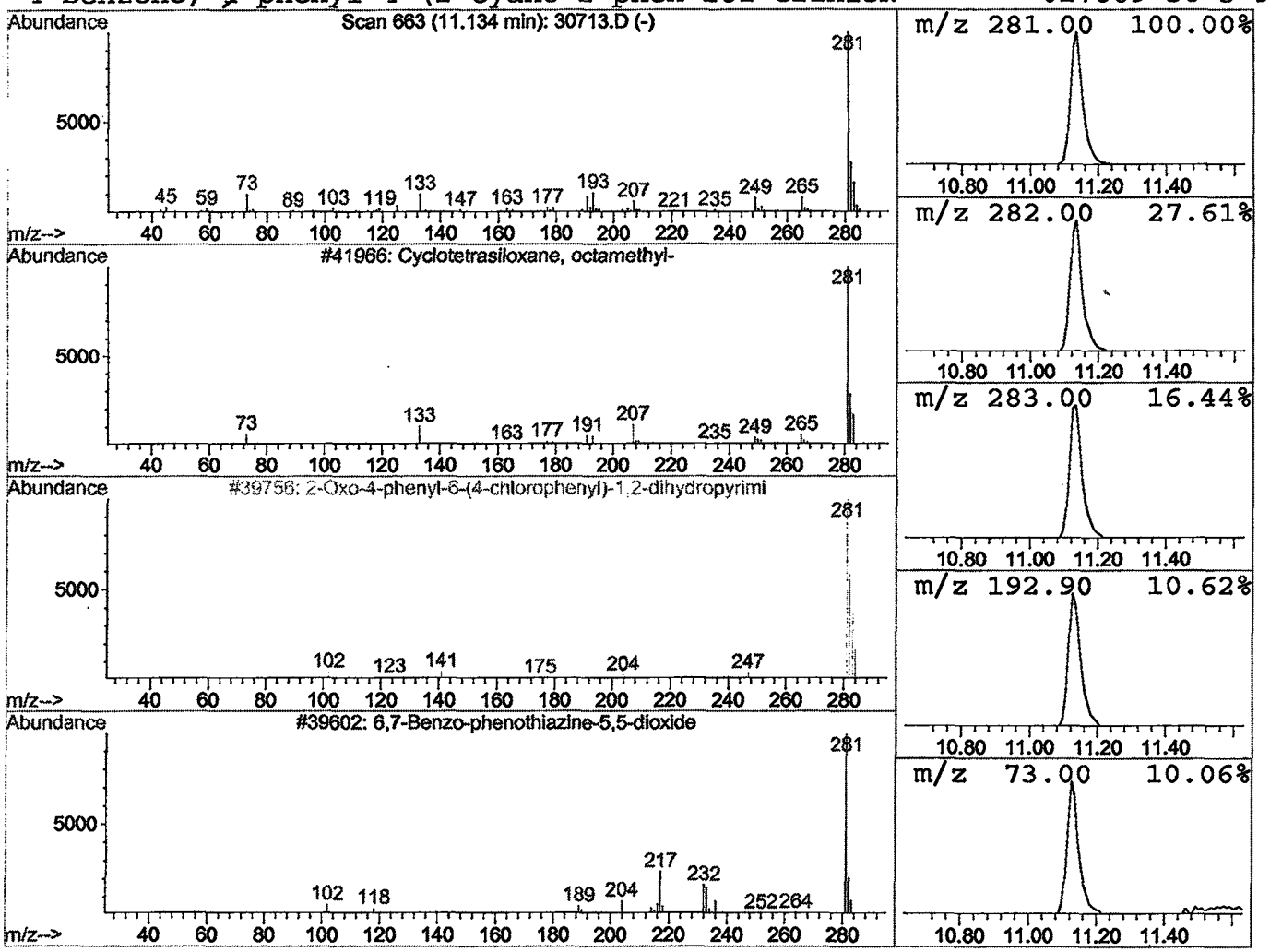
Quant Method : C:\HPCHEM\1\METHODS\GCMS1126.M (RTE Integrator)
Title : 209 625/TCLP calibration table
Library : C:\DATABASE\NBS75K.L

****

Peak Number 2 Cyclotetrasiloxane, octamethyl Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.13	1.27 ug/l	363659	1,4-Dichlorobenzene-d4	11.68

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclotetrasiloxane, octamethyl-	296	C8H24O4Si4	000556-67-2	86
2	2-Oxo-4-phenyl-6-(4-chlorophenyl)-1	282	C16H11ClN2O	000000-00-0	9
3	6,7-Benzo-phenothiazine-5,5-dioxide	281	C16H11NO2S	000000-00-0	9
4	Benzene, 1-phenyl-4-(2-cyano-2-phen	281	C21H15N	027869-56-3	9



Library Search Compound Report

Data File : C:\HPCHEM\1\DATA\DEC3101\30713.D Vial: 28
Acq On : 1 Jan 2002 3:07 pm Operator: 209 GALOS
Sample : gcms scan 12/19 Inst : GC/MS 209
Misc : Multiplr: 1.00
MS Integration Params: LSCINT.P

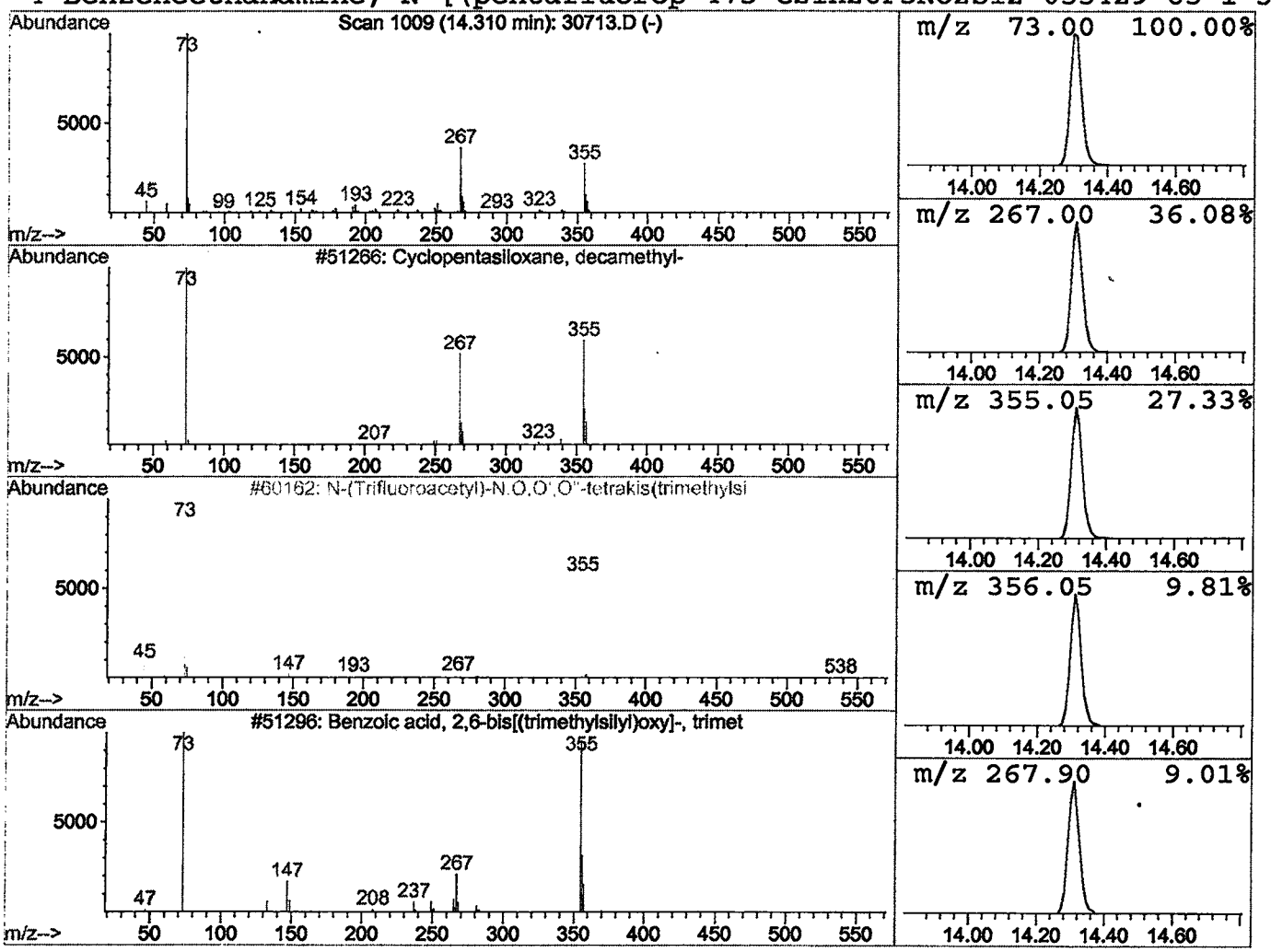
Quant Method : C:\HPCHEM\1\METHODS\GCMS1126.M (RTE Integrator)
Title : 209 625/TCLP calibration table
Library : C:\DATABASE\NBS75K.L

****

Peak Number 3 Cyclopentasiloxane, decamethyl Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.31	2.32 ug/l	816347	Naphthalene-d8	15.28

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclopentasiloxane, decamethyl-	370	C10H30O5Si5	000541-02-6	72
2	N-(Trifluoroacetyl)-N,O,O',O''-tetr	553	C22H42F3NO4Si4	000000-00-0	43
3	Benzoic acid, 2,6-bis[(trimethylsil	370	C16H30O4Si3	003782-85-2	40
4	Benzeneethanamine, N-[(pentafluorop	475	C21H26F5NO2Si2	055429-85-1	38



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Acq On : 1 Jan 2002 3:07 pm Operator: 209 GALOS
Sample : gcms scan 12/19 Inst : GC/MS 209
Misc : Multiplr: 1.00
MS Integration Params: LSCINT.P

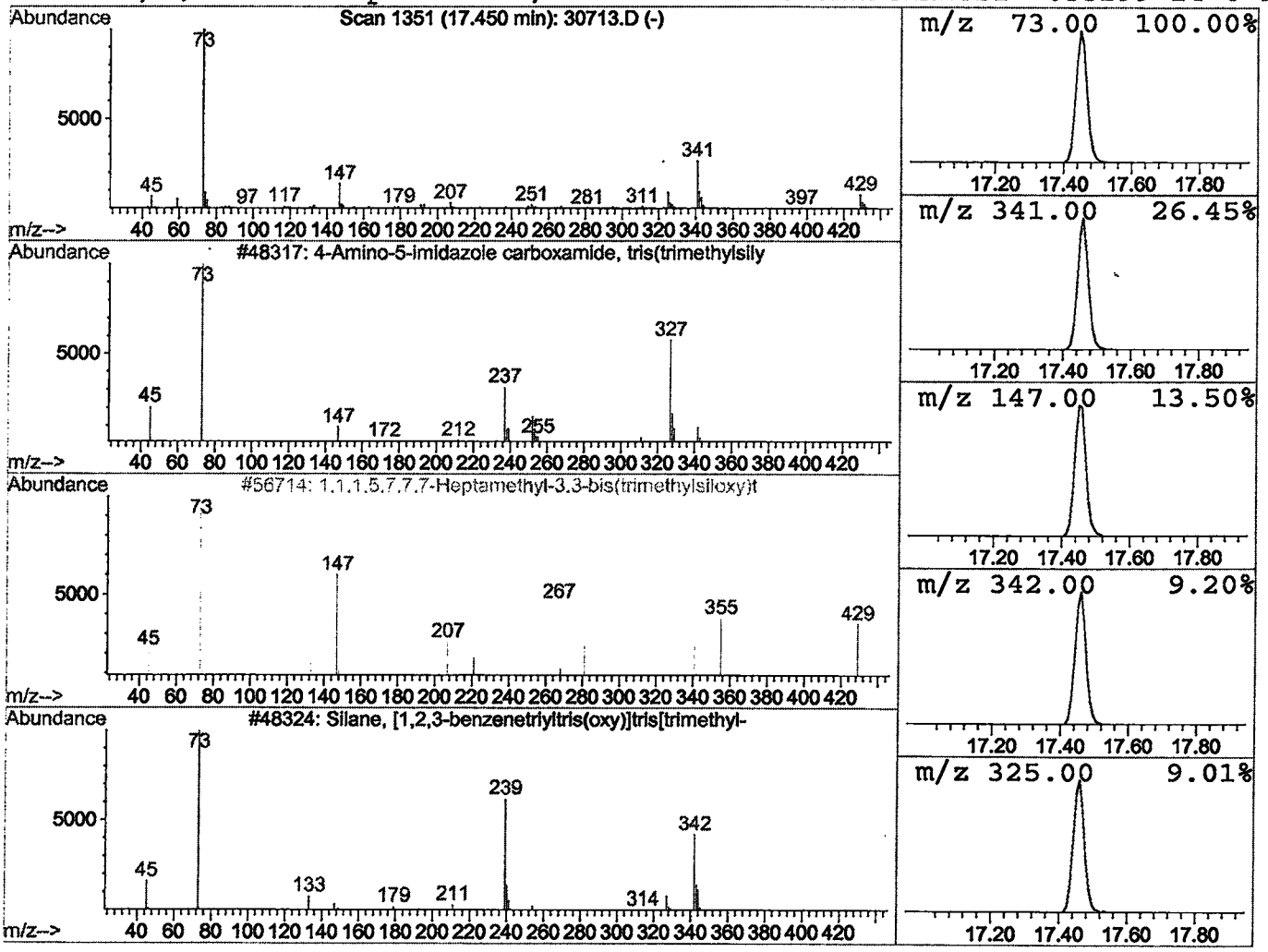
Quant Method : C:\HPCHEM\1\METHODS\GCMS1126.M (RTE Integrator)
Title : 209 625/TCLP calibration table
Library : C:\DATABASE\NBS75K.L

****

Peak Number 4 4-Amino-5-imidazole carboxamid Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.45	3.31 ug/l	1161250	Naphthalene-d8	15.28

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	4-Amino-5-imidazole carboxamide, tr	342	C13H30N4OSi3	000000-00-0	32
2	1,1,1,5,7,7,7-Heptamethyl-3,3-bis(t	444	C13H40O5Si6	038147-00-1	25
3	Silane, [1,2,3-benzenetriyltris(oxy	342	C15H30O3Si3	017864-23-2	10
4	2H-1,4-Benzodiazepin-2-one, 7-chlor	342	C18H19ClN2OSi	055299-24-6	9



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Acq On : 1 Jan 2002 3:07 pm
Sample : gcms scan 12/19
Misc :
MS Integration Params: LSCINT.P

Vial: 28
Operator: 209 GALOS
Inst : GC/MS 209
Multiplr: 1.00

Quant Method : C:\HPCHEM\1\METHODS\GCMS1126.M (RTE Integrator)
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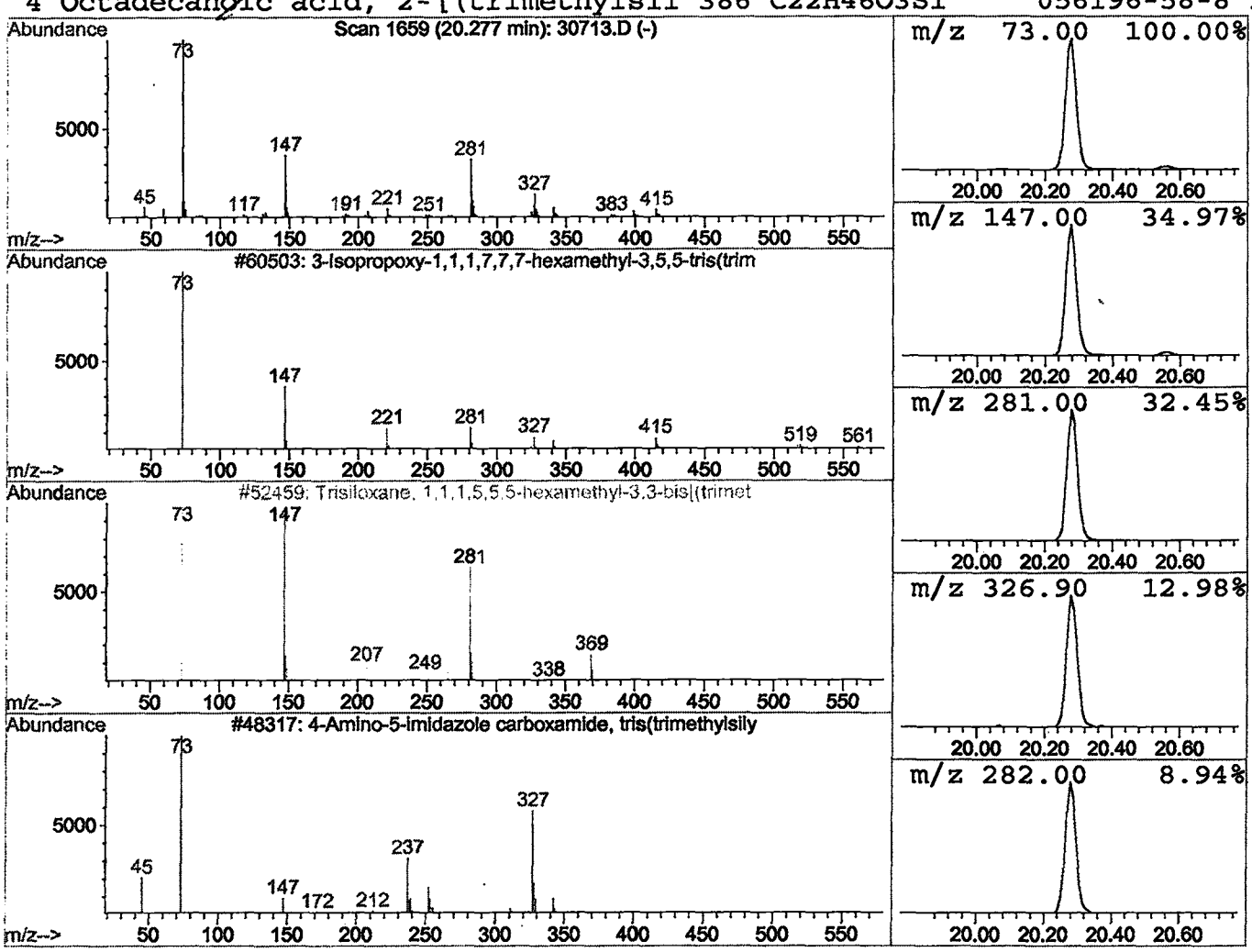
****

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.28	1.93 ug/l	762328	Acenaphthene-d10	20.56

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	3-Isopropoxy-1,1,1,7,7,7-hexamethyl	576	C18H52O7Si7	071579-69-6	45
2	Trisiloxane, 1,1,1,5,5,5-hexamethyl	384	C12H36O4Si5	003555-47-3	43
3	4-Amino-5-imidazole carboxamide, tr	342	C13H30N4OSi3	000000-00-0	16
4	Octadecanoic acid, 2-[(trimethylsil	386	C22H46O3Si	056196-58-8	10



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Sample : gcms scan 12/19
Misc :
MS Integration Params: LSCINT.P

Vial: 28
Operator: 209 GALOS
Inst : GC/MS 209
Multiplr: 1.00

Quant Method : C:\HPCHEM\1\METHODS\GCMS1126.M (RTE Integrator)
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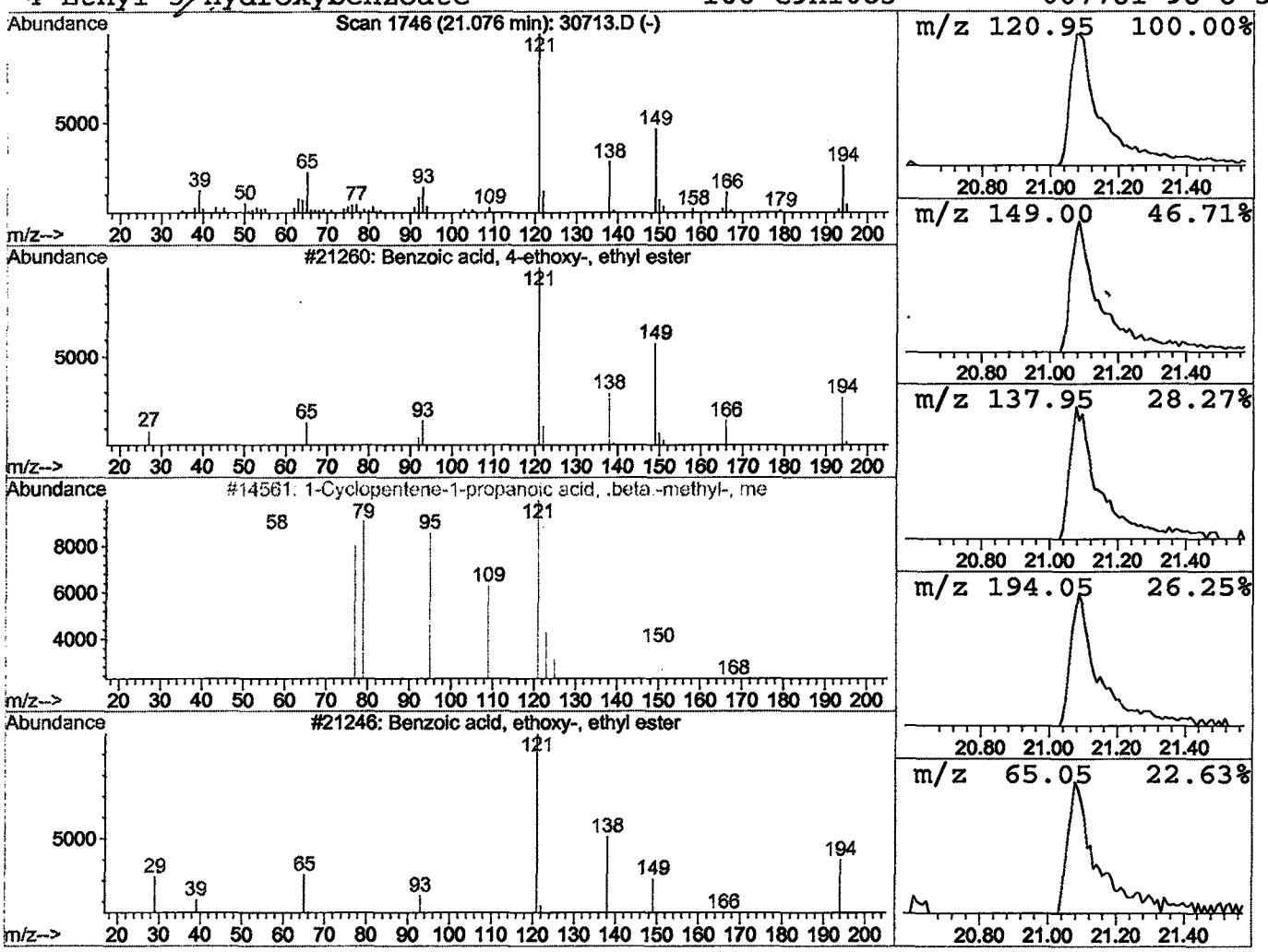
Library : C:\DATABASE\NBS75K.L

****

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.08	0.63 ug/l	247350	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	95
2	1-Cyclopentene-1-propanoic acid, .b	168	C10H16O2	061833-30-5	42
3	Benzoic acid, ethoxy-, ethyl ester	194	C11H14O3	075333-22-1	38
4	Ethyl 3-hydroxybenzoate	166	C9H10O3	007781-98-8	38



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Sample : gcms scan 12/19
Misc :
MS Integration Params: LSCINT.P

Vial: 28
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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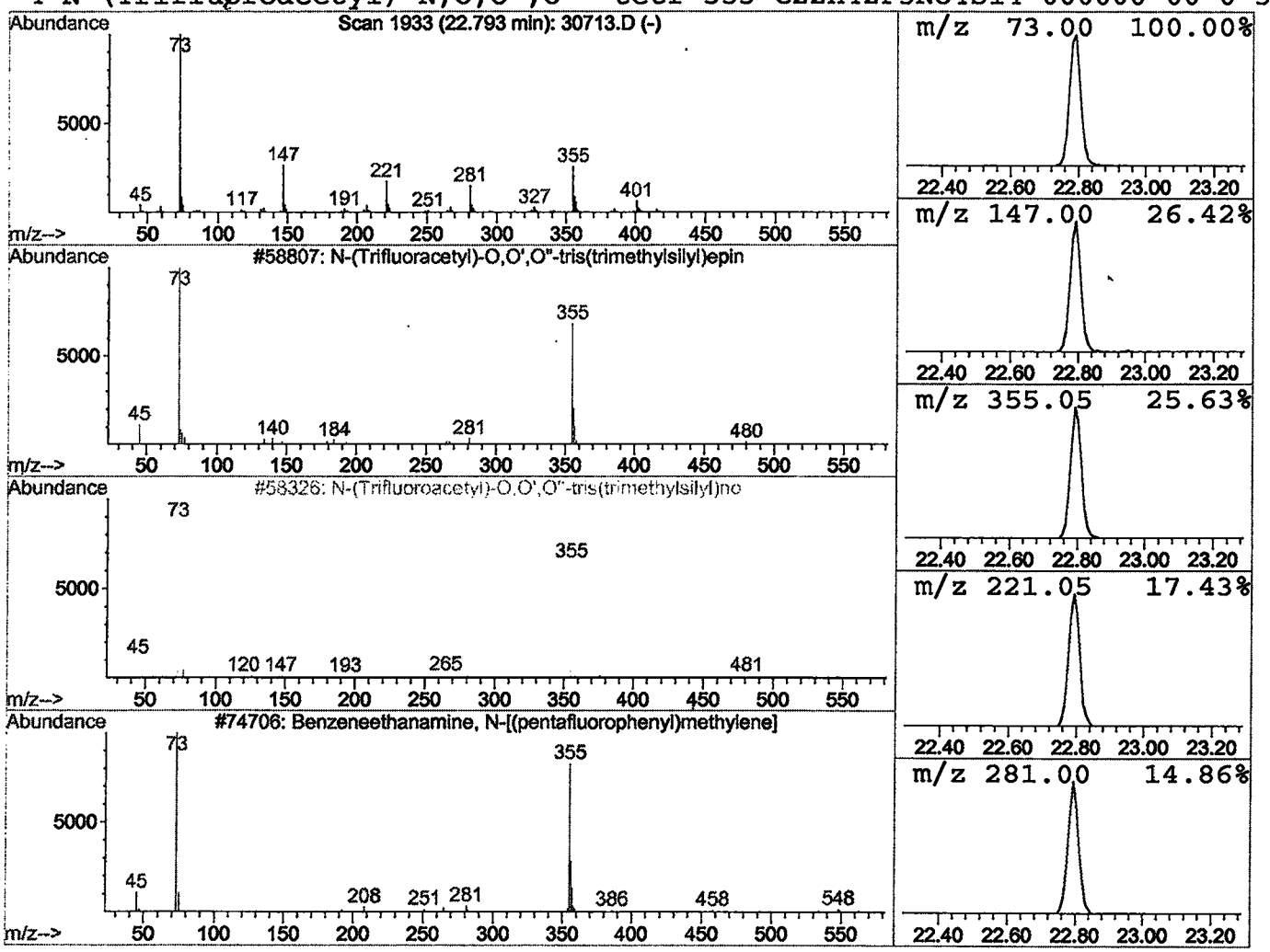
Peak Number 7 N-(Trifluoracetyl)-O,O',O"-tri Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.79	1.18 ug/l	464199	Phenanthrene-d10	24.99

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	N-(Trifluoracetyl)-O,O',O"-tris(tri	495	C20H36F3NO4Si3	000000-00-0	37
2	N-(Trifluoroacetyl)-O,O',O''-tris(t	481	C19H34F3NO4Si3	000000-00-0	37
3	Benzeneethanamine, N-[(pentafluorop	563	C24H34F5NO3Si3	055429-13-5	37
4	N-(Trifluoroacetyl)-N,O,O',O''-tetr	553	C22H42F3NO4Si4	000000-00-0	35



Library Search Compound Report

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Acq On : 1 Jan 2002 3:07 pm
Sample : gcms scan 12/19
Misc :
MS Integration Params: LSCINT.P

Vial: 28
Operator: 209 GALOS
Inst : GC/MS 209
Multiplr: 1.00

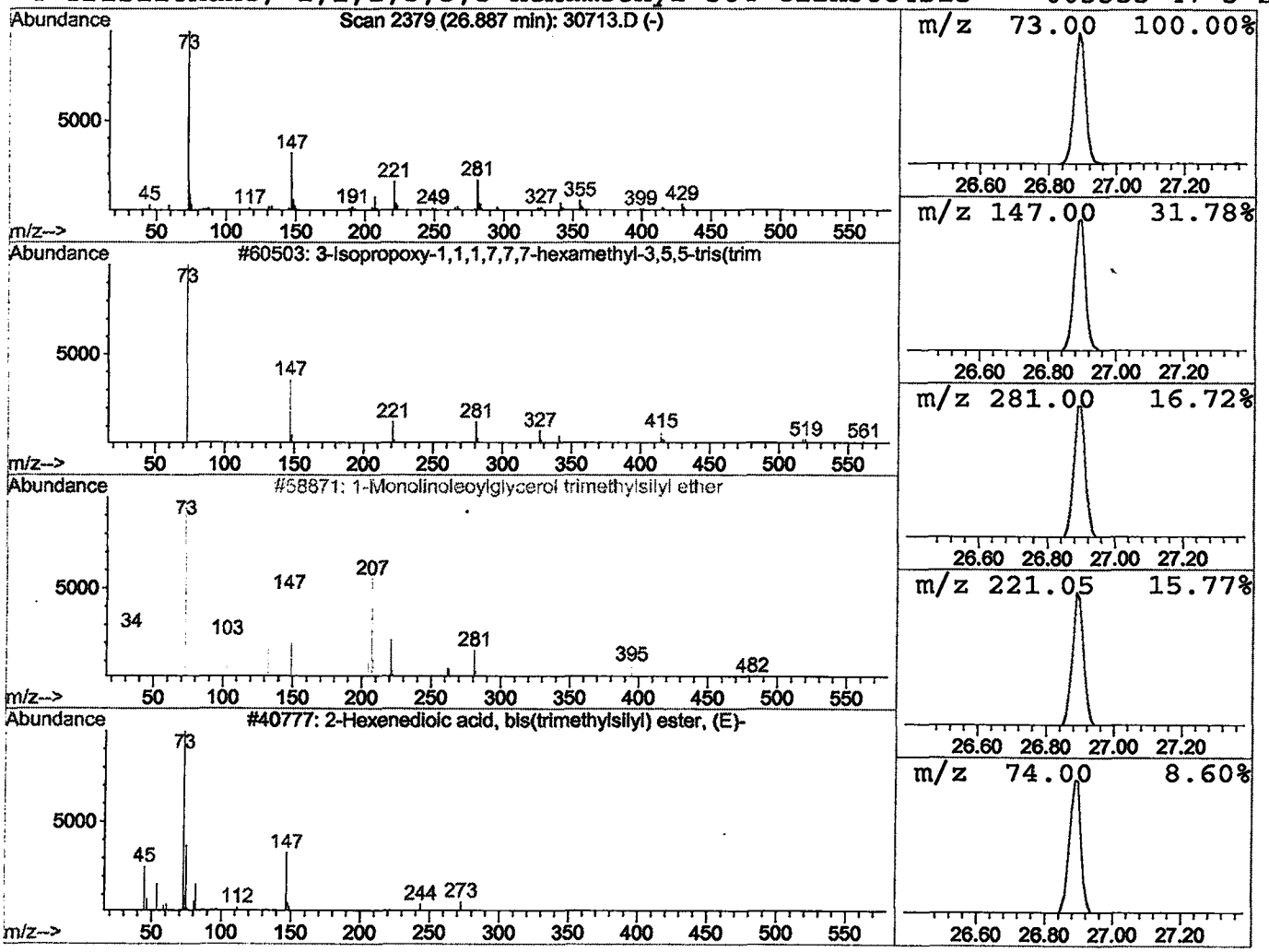
Quant Method : C:\HPCHEM\1\METHODS\GCMS1126.M (RTE Integrator)
Title : 209 625/TCLP calibration table
Library : C:\DATABASE\NBS75K.L

****

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	0.51 ug/l	201540	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	50
2	1-Monolinoleoylglycerol trimethylsi	498	C27H54O4Si2	054284-45-6	38
3	2-Hexenedioic acid, bis(trimethylsi	288	C12H24O4Si2	055494-10-5	32
4	Trisiloxane, 1,1,1,5,5,5-hexamethyl	384	C12H36O4Si5	003555-47-3	25



Library Search Compound Report

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Sample : gcms scan 12/19
Misc :
MS Integration Params: LSCINT.P

Vial: 28
Operator: 209 GALOS
Inst : GC/MS 209
Multiplr: 1.00

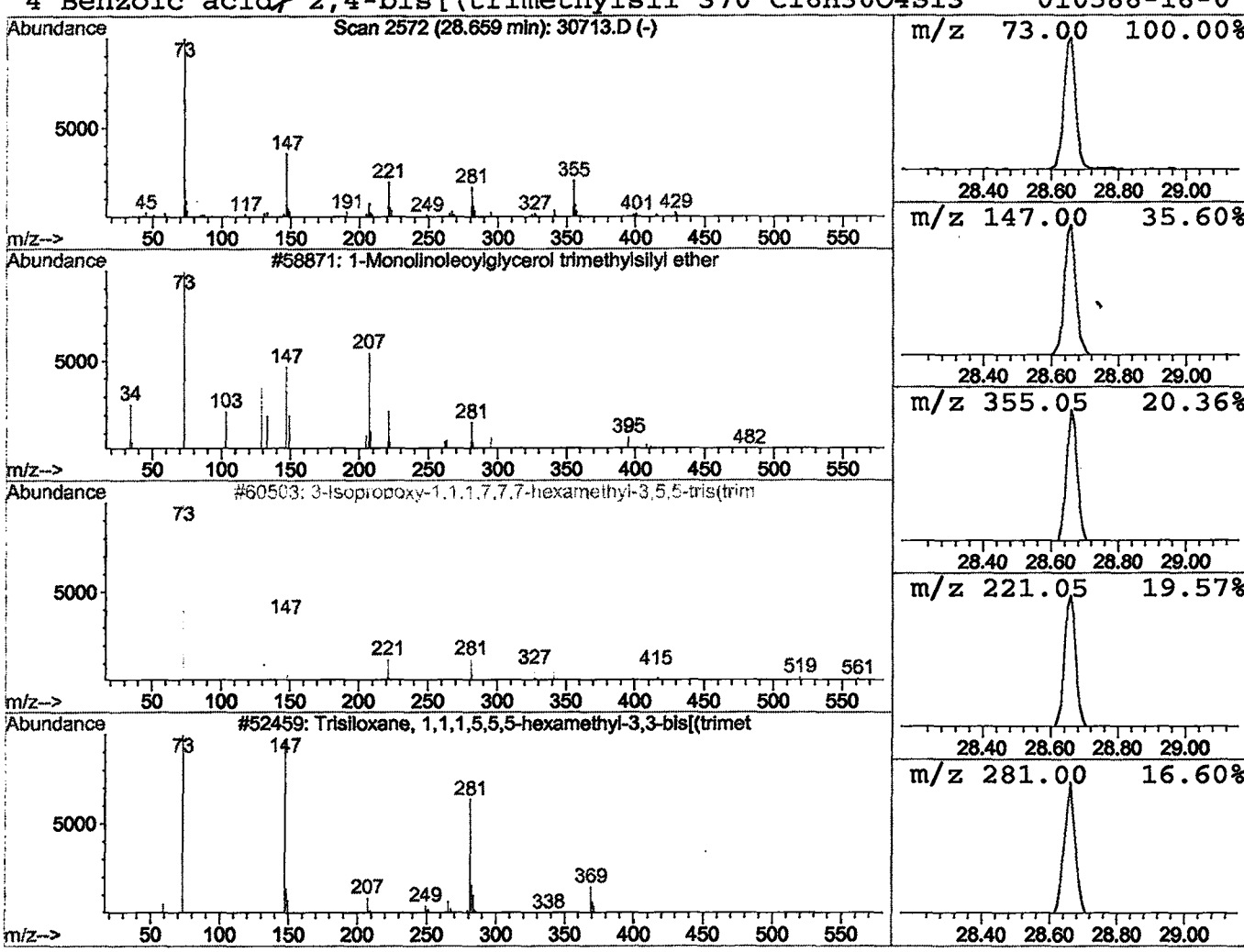
Quant Method : C:\HPCHEM\1\METHODS\GCMS1126.M (RTE Integrator)
Title : 209 625/TCLP calibration table
Library : C:\DATABASE\NBS75K.L

****

Peak Number 9 1-Monolinoleoylglycerol trimet Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
28.66	0.36 ug/l	142441	Phenanthrene-d10	24.99

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Monolinoleoylglycerol trimethylsi	498	C27H54O4Si2	054284-45-6	33
2	3-Isopropoxy-1,1,1,7,7,7-hexamethyl	576	C18H52O7Si7	071579-69-6	25
3	Trisiloxane, 1,1,1,5,5,5-hexamethyl	384	C12H36O4Si5	003555-47-3	25
4	Benzoic acid, 2,4-bis[(trimethylsil	370	C16H30O4Si3	010586-16-0	12



Library Search Compound Report

Data File : C:\HPCHEM\1\DATA\DEC3101\30713.D
Acq On : 1 Jan 2002 3:07 pm
Sample : gcms scan 12/19
Misc :
MS Integration Params: LSCINT.P

Vial: 28
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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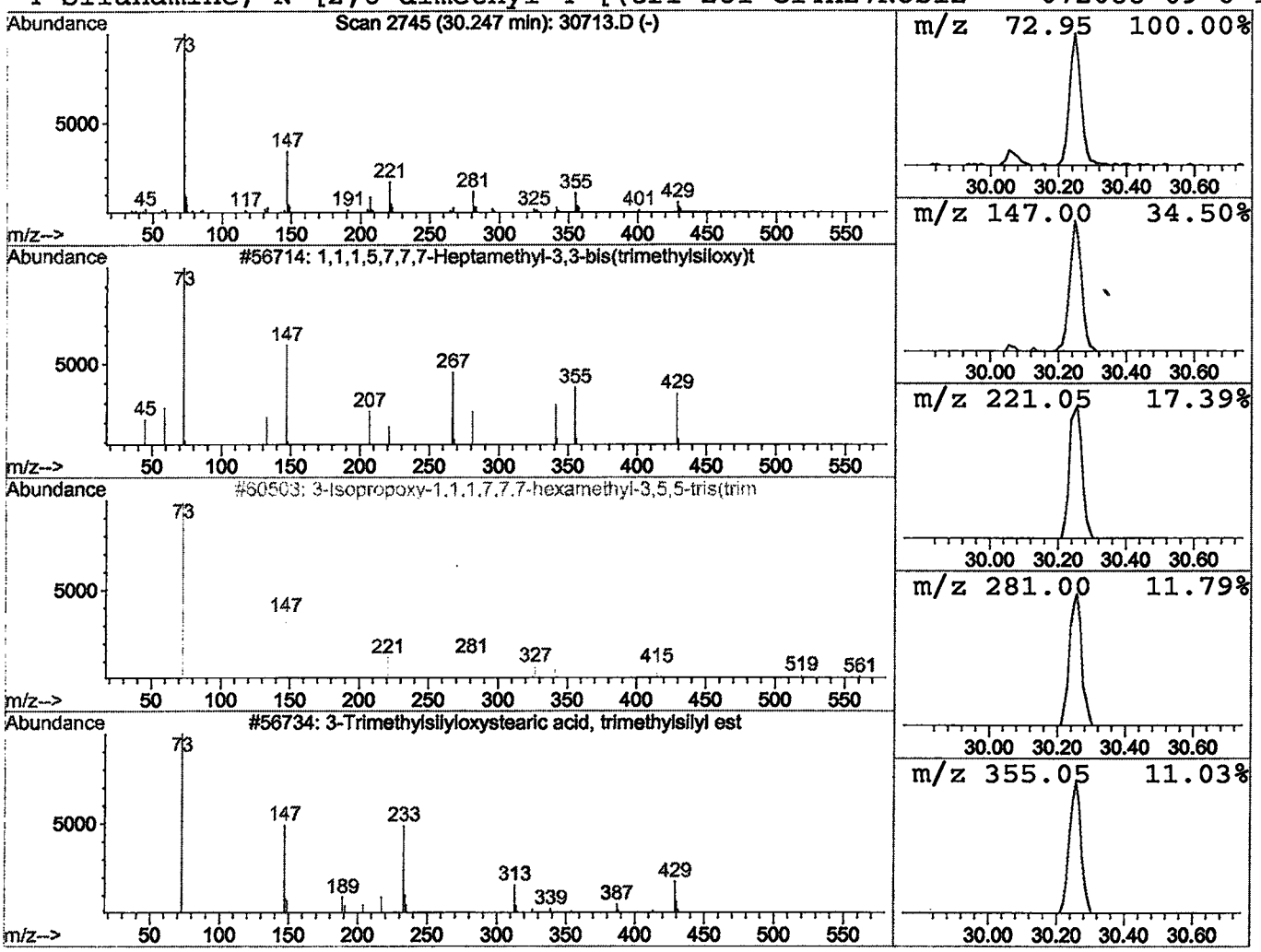
Peak Number 10 1,1,1,5,7,7,7-Heptamethyl-3,3- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
30.25	0.47 ug/l	123012	Chrysene-d12	33.02

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	34
2	3-Isopropoxy-1,1,1,7,7,7-hexamethyl	576	C18H52O7Si7	071579-69-6	32
3	3-Trimethylsilyloxystearic acid, tr	444	C24H52O3Si2	000000-00-0	17
4	Silanamine, N-[2,6-dimethyl-4-[(tri	281	C14H27NOSi2	072088-09-6	16



Library Search Compound Report

Data File : C:\HPCHEM\1\DATA\DEC3101\30713.D
Acq On : 1 Jan 2002 3:07 pm
Sample : gcms scan 12/19
Misc :
MS Integration Params: LSCINT.P

Vial: 28
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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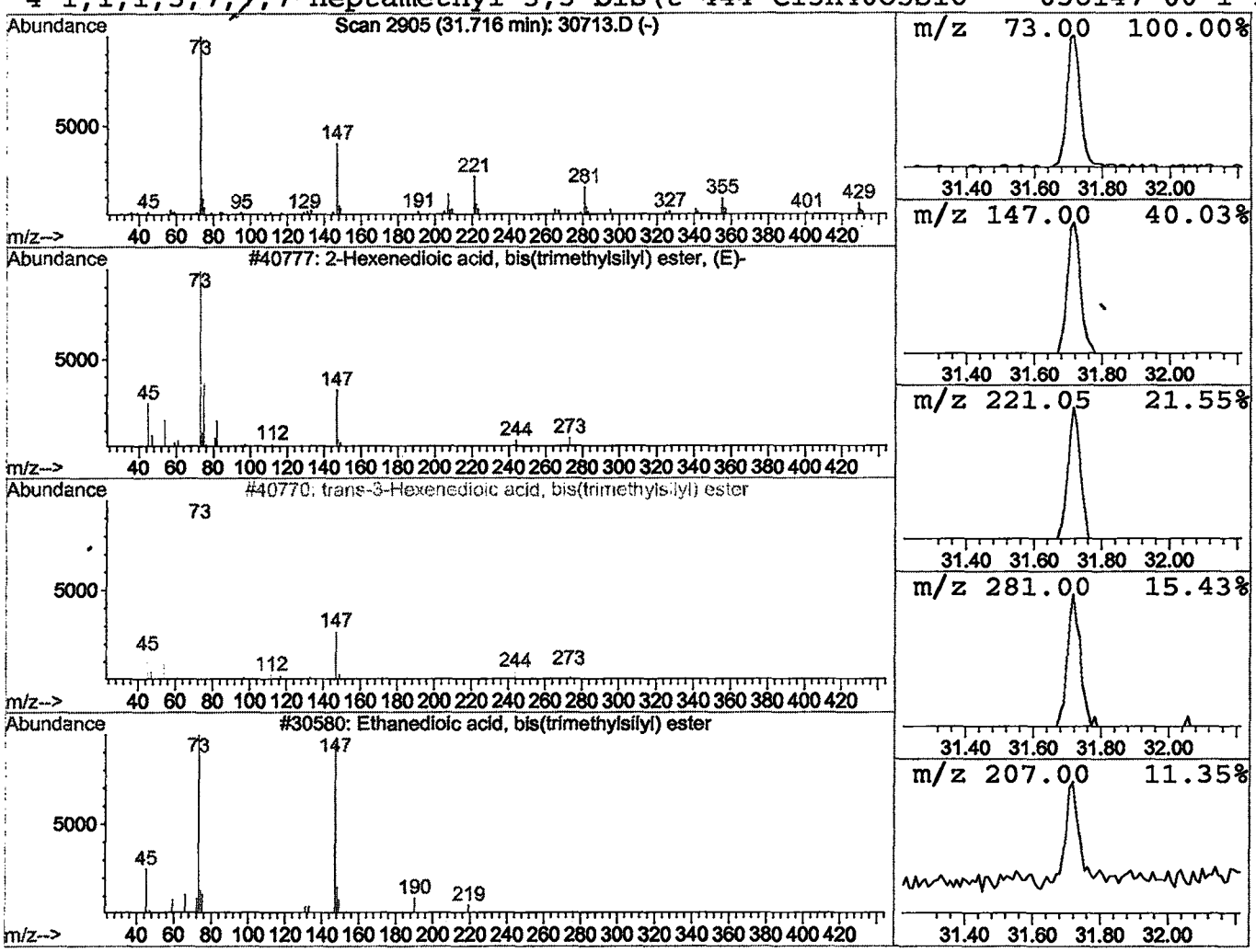
Peak Number 11 2-Hexenedioic acid, bis(trimet Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
31.72	0.39 ug/l	102071	Chrysene-d12	33.02

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Hexenedioic acid, bis(trimethylsi	288	C12H24O4Si2	055494-10-5	23
2	trans-3-Hexenedioic acid, bis(trime	288	C12H24O4Si2	000000-00-0	23
3	Ethanedioic acid, bis(trimethylsily	234	C8H18O4Si2	018294-04-7	23
4	1,1,1,5,7,7,7-Heptamethyl-3,3-bis(t	444	C13H40O5Si6	038147-00-1	13



Tentatively Identified Compound (LSC) summary

Operator ID: 209 GALOS Date Acquired: 1 Jan 2002 3:07 pm
Data File: C:\HPCHEM\1\DATA\DEC3101\30713.D
Name: gcms scan 12/19
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
3-Pentanol, 3-methyl	7.62	3.9	ug/l	1118120	ISTD01	11.68	5713710	20.0
Cyclotetrasiloxane,	11.13	1.3	ug/l	363659	ISTD01	11.68	5713710	20.0
Cyclopentasiloxane,	14.31	2.3	ug/l	816347	ISTD02	15.28	7027210	20.0
4-Amino-5-imidazole	17.45	3.3	ug/l	1161250	ISTD02	15.28	7027210	20.0
3-Isopropoxy-1,1,1,7	20.28	1.9	ug/l	762328	ISTD03	20.56	7912600	20.0
Benzoic acid, 4-etho	21.08	0.6	ug/l	247350	ISTD03	20.56	7912600	20.0
N-(Trifluoracetyl)-O	22.79	1.2	ug/l	464199	ISTD04	24.99	7840910	20.0
3-Isopropoxy-1,1,1,7	26.89	0.5	ug/l	201540	ISTD04	24.99	7840910	20.0
1-Monolinoleoylglyce	28.66	0.4	ug/l	142441	ISTD04	24.99	7840910	20.0
1,1,1,5,7,7,7-Heptam	30.25	0.5	ug/l	123012	ISTD05	33.02	5195460	20.0
2-Hexenedioic acid,	31.72	0.4	ug/l	102071	ISTD05	33.02	5195460	20.0

30713.D GCMS1126.M Wed Jan 23 08:47:08 2002

column bleed