User: Vinci, David L
Westchester Dept. of Labs & Research
Date: 01/10/2002 Time: 08:42:06

Sample: AD30426
Analysis: $GCMS GCMS Scan For Unknowns
Units: ug
Started: 1/10/02 08:37 Ended: 1/10/02 08:37 Analyst: DLV
Result source: manual entry
Page: 1

	Component Name	Viol	Result	Qualifier	MDL
1	Other unknown compounds		.		
2	Surr_2,4-DDD		89		5.0

Comments for Sample AD30426 - Analysis $GCMS

Westchester Dept. of Labs & Research
User: Vinci, David L
Date: 01-10-2002 Time: 08:42:09

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\30426.D Vial: 23
Acq On : 28 Dec 2001 11:28 am Operator: 209 GALOS
Sample : GC/MS SCAN 12/13 Inst : GC/MS 209
Misc : Multiplr: 1.00
MS Integration Params: GOOFY.P
Quant Time: Dec 31 12:19 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 14 12:19:30 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	1079887	20.00	ug/l	-0.02
10) Naphthalene-d8	15.28	136	4159467	20.00	ug/l	-0.03
17) Acenaphthene-d10	20.56	164	2115045	20.00	ug/l	-0.02
29) Phenanthrene-d10	24.98	188	3168694m	20.00	ug/l	-0.01
38) Chrysene-d12	33.02	240	1964485	20.00	ug/l	-0.01
44) Perylene-d12	37.12	264	1146695	20.00	ug/l	0.00
System Monitoring Compounds						
37) 2,4-DDD	30.07	235	204607	4.46	ug/mL	0.00
Spiked Amount 5.000			Recovery = 89.20%			
Target Compounds						Qvalue
2) N-nitroso-dimethyl amine	5.11	74	4450	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	13.26	77	198	N.D.		
12) Isophorone	13.50	82	333	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	1357	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	989	N.D.		
21) 2,6-Dinitrotoluene	20.27	165	525	N.D.		
22) Acenaphthylene	0.00	152	0	N.D.		
23) Acenaphthene	20.57	153	142	N.D.		
24) 2,4-Dinitrotoluene	21.33	165	277	N.D.		
25) Diethylphthalate	22.13	149	2534	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\30426.D Vial: 23
Acq On : 28 Dec 2001 11:28 am Operator: 209 GALOS
Sample : GC/MS SCAN 12/13 Inst : GC/MS 209
Misc : Multiplr: 1.00
MS Integration Params: GOOFY.P
Quant Time: Dec 31 12:19 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 14 12:19:30 2001
Response via : Initial Calibration
DataAcq Meth : GCMS1126

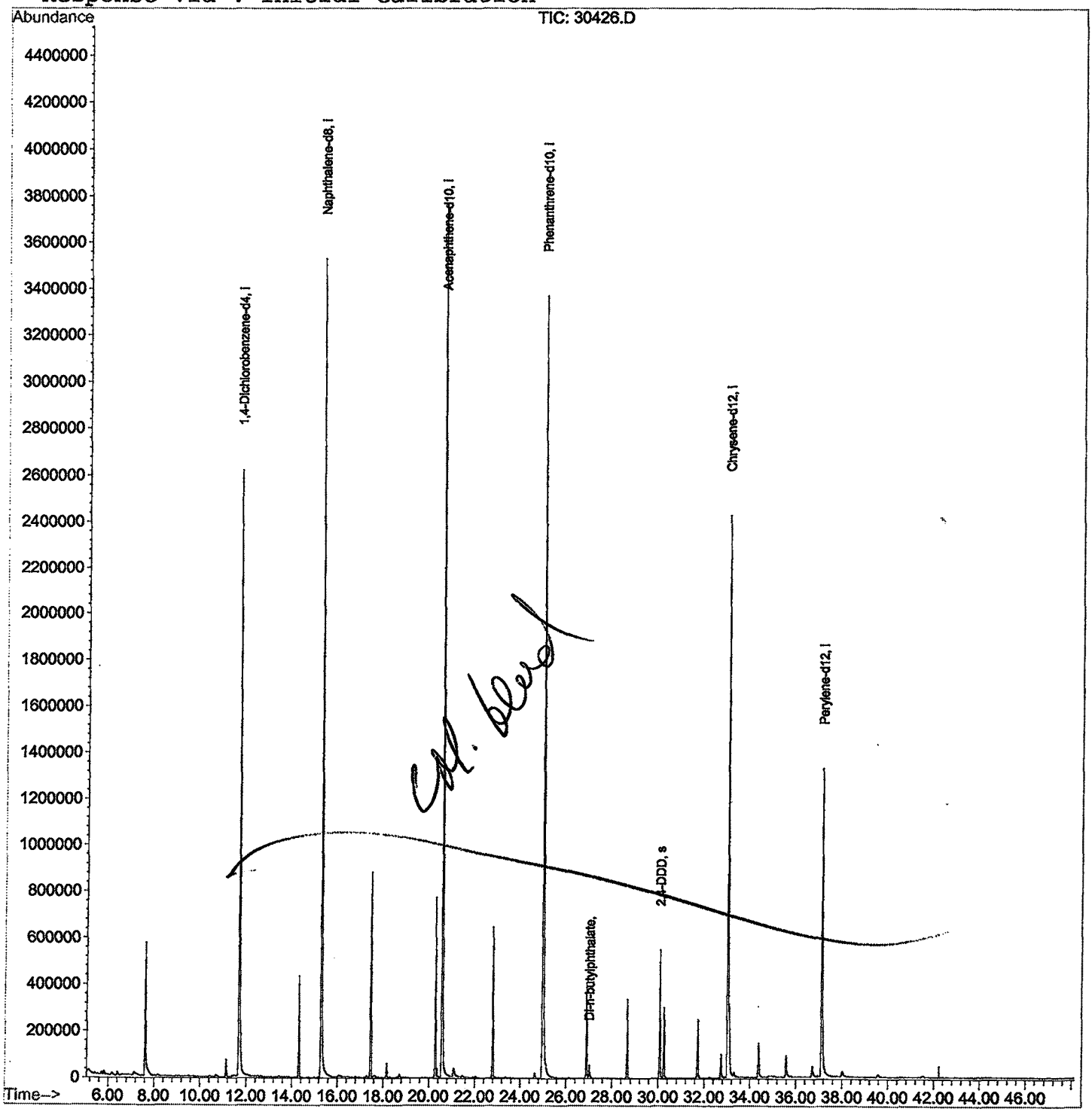
	Compound	R.T.	QIon	Response	Conc Unit	Qvalue
30)	N-Nitrosodiphenylamine	0.00	168	0	N.D.	
31)	4-Bromophenyl-phenylether	0.00	248	0	N.D.	
32)	Hexachlorobenzene	0.00	284	0	N.D.	
33)	Phenanthrene	24.98	178	3176	N.D.	
34)	Anthracene	25.18	178	472	N.D.	
35)	Di-n-butylphthalate	27.02	149	92618	0.41 ug/l	
36)	Fluoranthene	0.00	202	0	N.D.	
39)	Pyrene	29.47	202	763	N.D.	
40)	Butylbenzylphthalate	31.53	149	1264	N.D.	
41)	Benzo(a)anthracene	33.09	228	7411	N.D.	
42)	Bis(2-ethylhexyl)phthalate	33.31	149	13538	N.D.	
43)	Chrysene	33.09	228	7411	N.D.	
45)	Di-n-Octylphthalate	35.05	149	6310	N.D.	
46)	Benzo(b)fluoranthene	36.09	252	2104	N.D.	
47)	Benzo(k)fluoranthene	36.09	252	2367	N.D.	
48)	Benzo(a)pyrene	36.95	252	1515	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\30426.D Vial: 23
Acq On : 28 Dec 2001 11:28 am Operator: 209 GALOS
Sample : GC/MS SCAN 12/13 Inst : GC/MS 209
Misc : Multiplr: 1.00
MS Integration Params: GOOFY.P
Quant Time: Dec 31 12:19 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 : C:\HPCHEM\1\DATA\DEC2701\30426.D Vial: 23
Acq On : 28 Dec 2001 11:28 am Operator: 209 GALOS
Sample : GC/MS SCAN 12/13 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

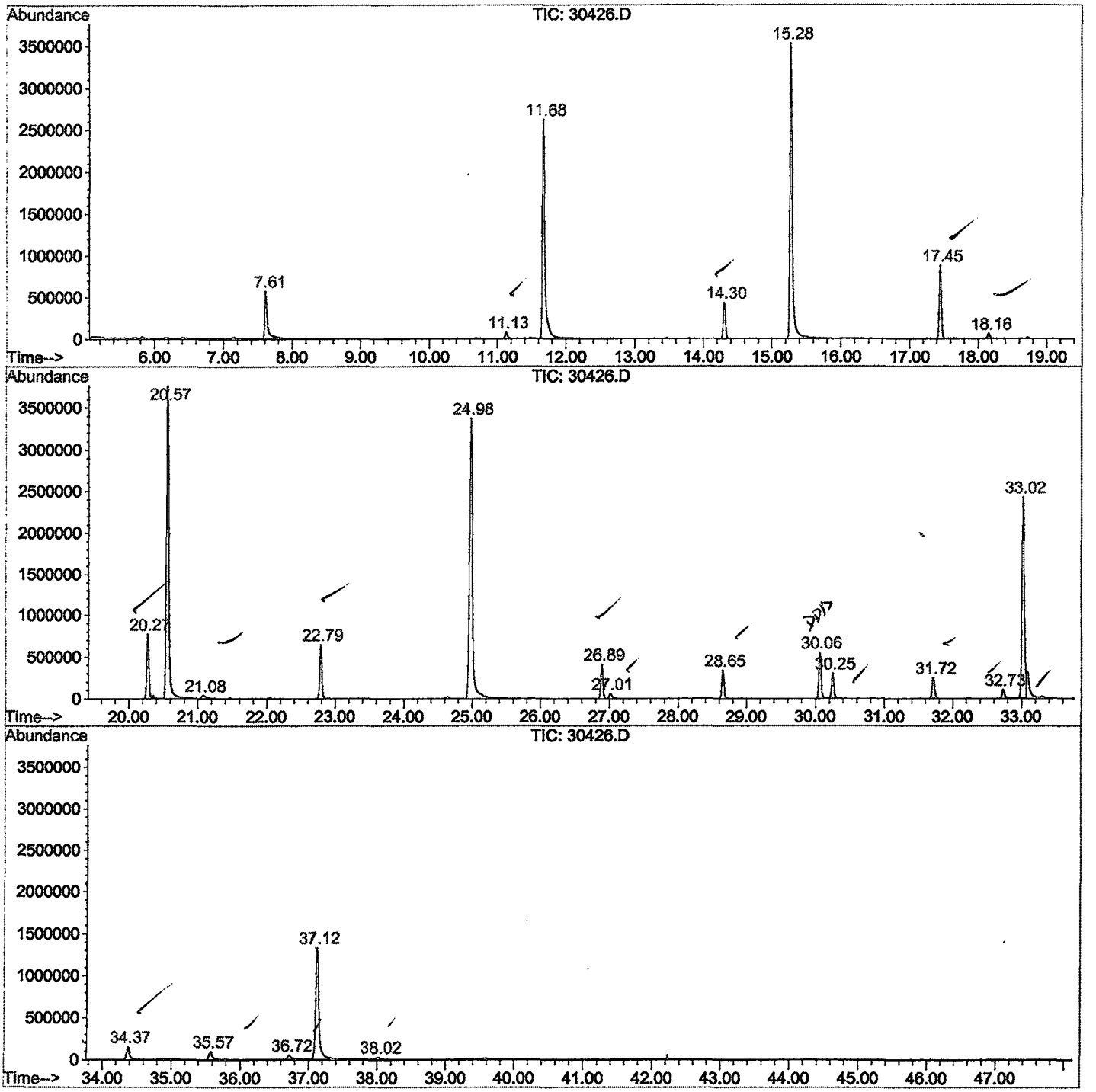
peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	peak area	peak % max.	% of total
1	7.612	276	280	310	rBV	569311	1480911	14.89%	2.556%
2	11.128	656	663	674	rVB	75580	174510	1.75%	0.301%
3	11.678	717	723	747	rBV	2616568	6639581	66.77%	11.461%
4	14.304	1001	1009	1017	rBV	433505	929424	9.35%	1.604%
5	15.277	1109	1115	1135	rBV	3526452	8359549	84.06%	14.430%
6	17.453	1344	1352	1362	rBV	881296	1870656	18.81%	3.229%
7	18.160	1424	1429	1440	rBV	62323	139029	1.40%	0.240%
8	20.271	1652	1659	1667	rBV	771723	1614362	16.23%	2.787%
9	20.565	1684	1691	1720	rBV	3768249	9309613	93.61%	16.070%
10	21.079	1741	1747	1759	rBV2	34441	154884	1.56%	0.267%
11	22.787	1927	1933	1949	rVB	646424	1344982	13.52%	2.322%
12	24.981	2163	2172	2202	rBV3	3370438	9944588	100.00%	17.166%
13	26.890	2373	2380	2387	rBV	406966	869295	8.74%	1.501%
14	27.010	2387	2393	2407	rVV	55191	163711	1.65%	0.283%
15	28.653	2565	2572	2581	rBV	337963	734867	7.39%	1.269%
16	30.057	2720	2725	2739	rBV	548178	1267323	12.74%	2.188%
17	30.250	2740	2746	2756	rVV	300320	694473	6.98%	1.199%
18	31.719	2899	2906	2915	rBV	249006	628540	6.32%	1.085%
19	32.729	3011	3016	3026	rBV	102845	274017	2.76%	0.473%
20	33.023	3041	3048	3053	rBV2	2422895	6041608	60.75%	10.429%
21	34.372	3188	3195	3212	rBV	148392	482977	4.86%	0.834%
22	35.575	3320	3326	3339	rBV	92433	315476	3.17%	0.545%
23	36.722	3444	3451	3459	rBV2	44441	161734	1.63%	0.279%
24	37.117	3486	3494	3516	rBV2	1323893	4231905	42.55%	7.305%
25	38.017	3584	3592	3603	rBV3	22349	103388	1.04%	0.178%

Sum of corrected areas: 57931403

30426.D GCMS1126.M Wed Jan 02 11:07:23 2002

LSC Report - Integrated Chromatogram

File : C:\HPCHEM\1\DATA\DEC2701\30426.D
Operator : 209 GALOS
Acquired : 28 Dec 2001 11:28 am using AcqMethod GCMS1126
Instrument : GC/MS 209
Sample Name: GC/MS SCAN 12/13
Misc Info :
Vial Number: 23
Quant File :GCMS1126.RES (RTE Integrator)



30426.D GCMS1126.M Wed Jan 02 11:07:30 2002

Library Search Compound Report

Data File : C:\HPCHEM\1\DATA\DEC2701\30426.D
Acq On : 28 Dec 2001 11:28 am
Sample : GC/MS SCAN 12/13
Misc :
MS Integration Params: LSCINT.P

Vial: 23
Operator: 209 GALOS
Inst : GC/MS 209
Multiplr: 1.00

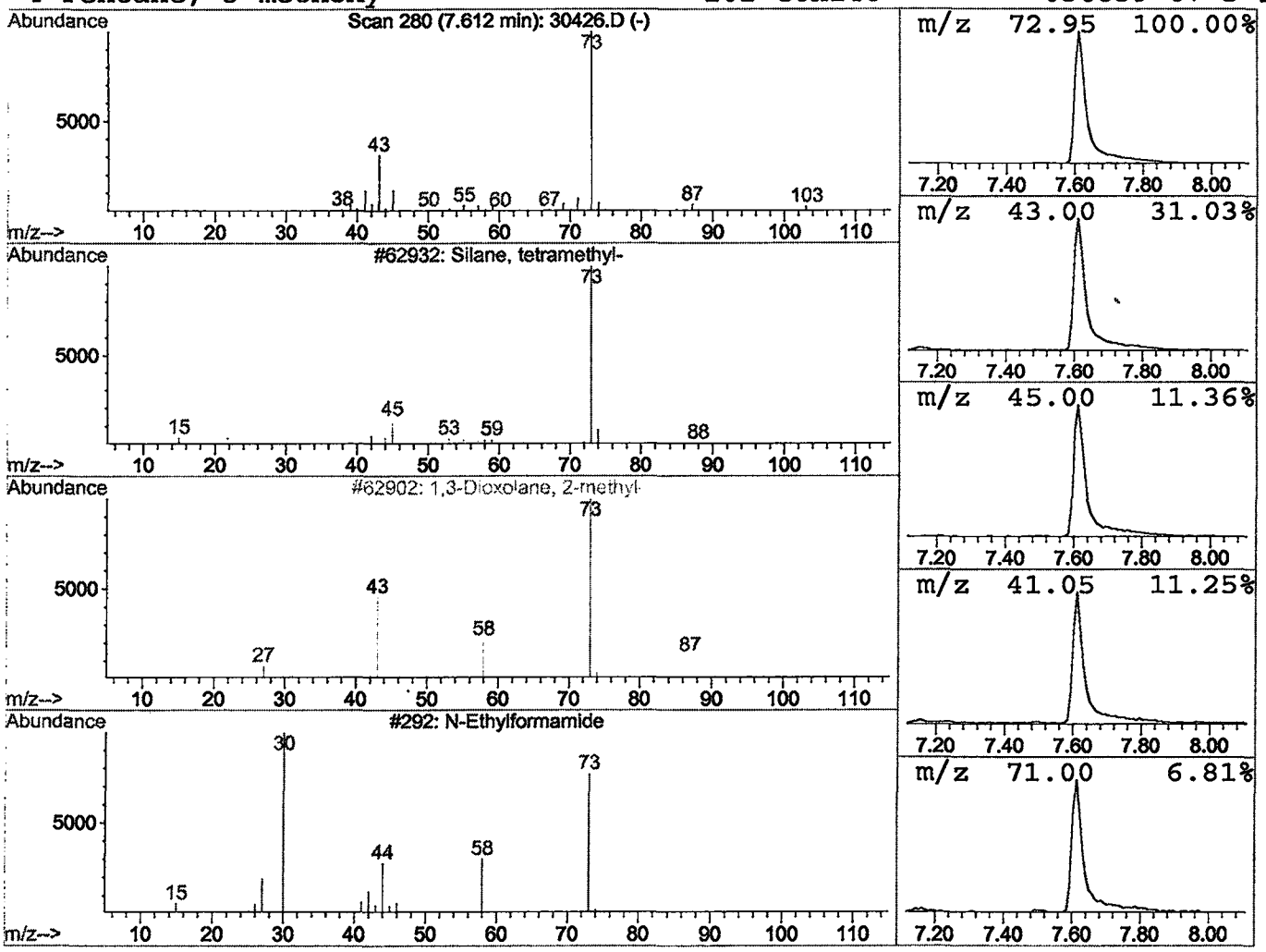
Quant Method : C:\HPCHEM\1\METHODS\GCMS1126.M (RTE Integrator)
Title : 209 625/TCLP calibration table
Library : C:\DATABASE\NBS75K.L

****

Peak Number 1 Silane, tetramethyl- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.61	4.46 ug/l	1480910	1,4-Dichlorobenzene-d4	11.68

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Silane, tetramethyl-	88	C4H12Si	000075-76-3	9
2	1,3-Dioxolane, 2-methyl-	88	C4H8O2	000497-26-7	9
3	N-Ethylformamide	73	C3H7NO	000627-45-2	9
4	Pentane, 3-methoxy-	102	C6H14O	036839-67-5	9



Library Search Compound Report

Data File : C:\HPCHEM\1\DATA\DEC2701\30426.D Vial: 23
Acq On : 28 Dec 2001 11:28 am Operator: 209 GALOS
Sample : GC/MS SCAN 12/13 Inst : GC/MS 209
Misc : Multiplr: 1.00
MS Integration Params: LSCINT.P

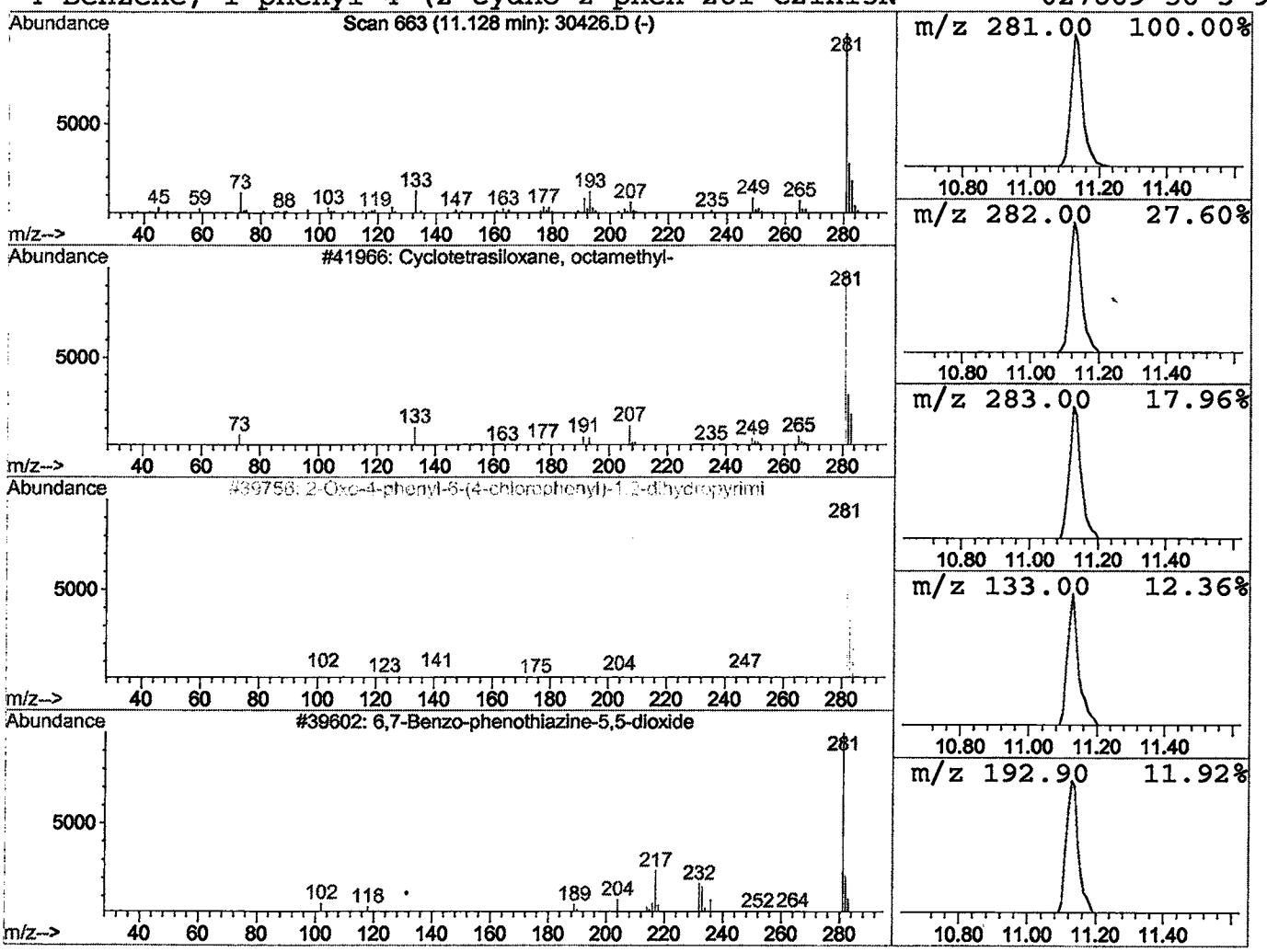
Quant Method : C:\HPCHEM\1\METHODS\GCMS1126.M (RTE Integrator)
Title : 209 625/TCLP calibration table
Library : C:\DATABASE\NBS75K.L

****

Peak Number 2 Cyclotetrasiloxane, octamethyl Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.13	0.53 ug/l	174510	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\DEC2701\30426.D
Acq On : 28 Dec 2001 11:28 am
Sample : GC/MS SCAN 12/13
Misc :
MS Integration Params: LSCINT.P

Vial: 23
Operator: 209 GALOS
Inst : GC/MS 209
Multiplr: 1.00

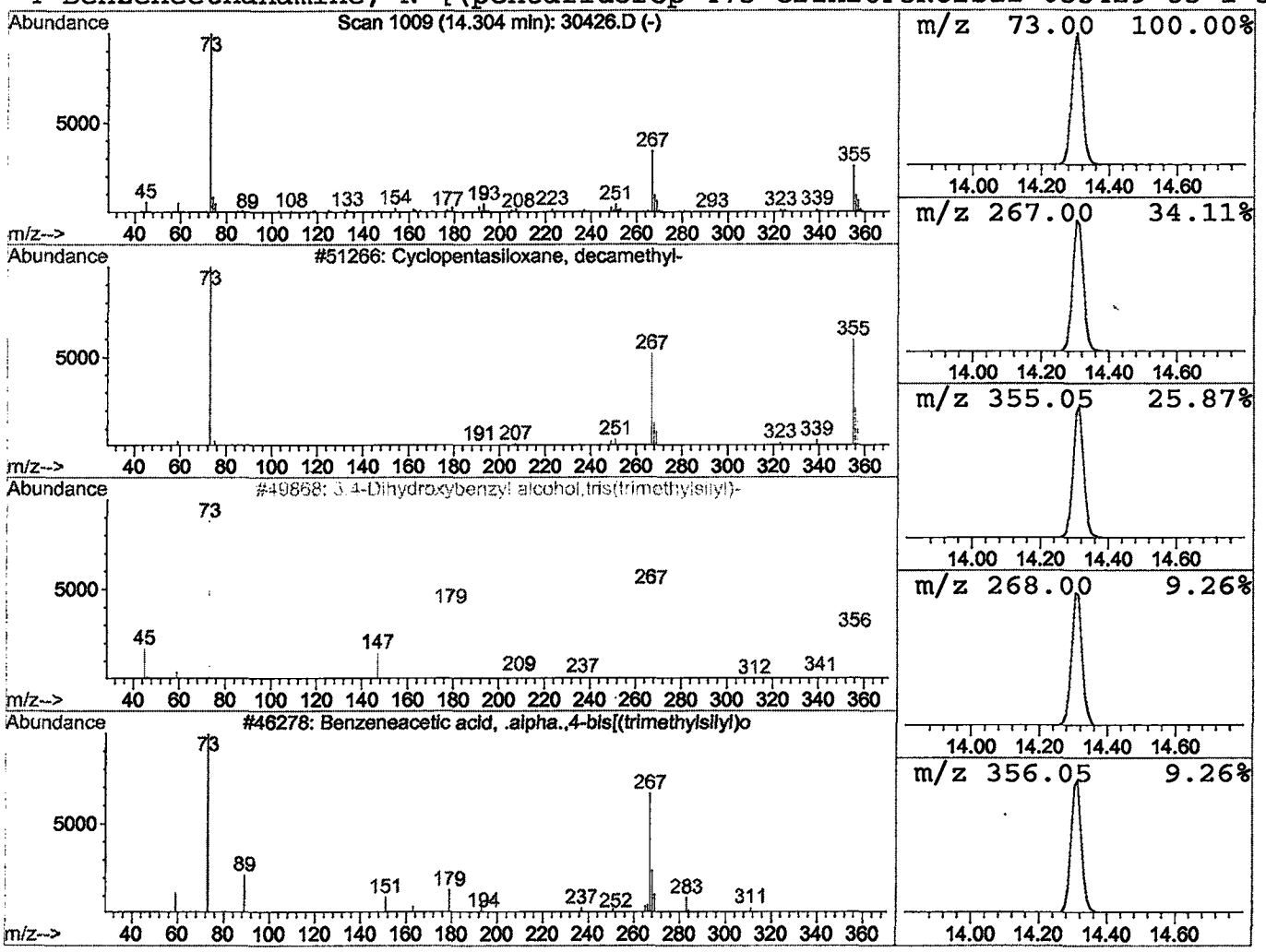
Quant Method : C:\HPCHEM\1\METHODS\GCMS1126.M (RTE Integrator)
Title : 209 625/TCLP calibration table
Library : C:\DATABASE\NBS75K.L

****

Peak Number 3 Cyclopentasiloxane, decamethyl Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.30	2.22 ug/l	929424	Naphthalene-d8	15.28

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclopentasiloxane, decamethyl-	370	C10H30O5Si5	000541-02-6	50
2	3,4-Dihydroxybenzyl alcohol,tris(tr	356	C16H32O3Si3	068595-79-9	38
3	Benzeneacetic acid, .alpha.,4-bis[(	326	C15H26O4Si2	055334-40-2	32
4	Benzeneethanamine, N-[(pentafluorop	475	C21H26F5NO2Si2	055429-85-1	32



Library Search Compound Report

Data File : C:\HPCHEM\1\DATA\DEC2701\30426.D
Acq On : 28 Dec 2001 11:28 am
Sample : GC/MS SCAN 12/13
Misc :
MS Integration Params: LSCINT.P

Vial: 23
Operator: 209 GALOS
Inst : GC/MS 209
Multiplr: 1.00

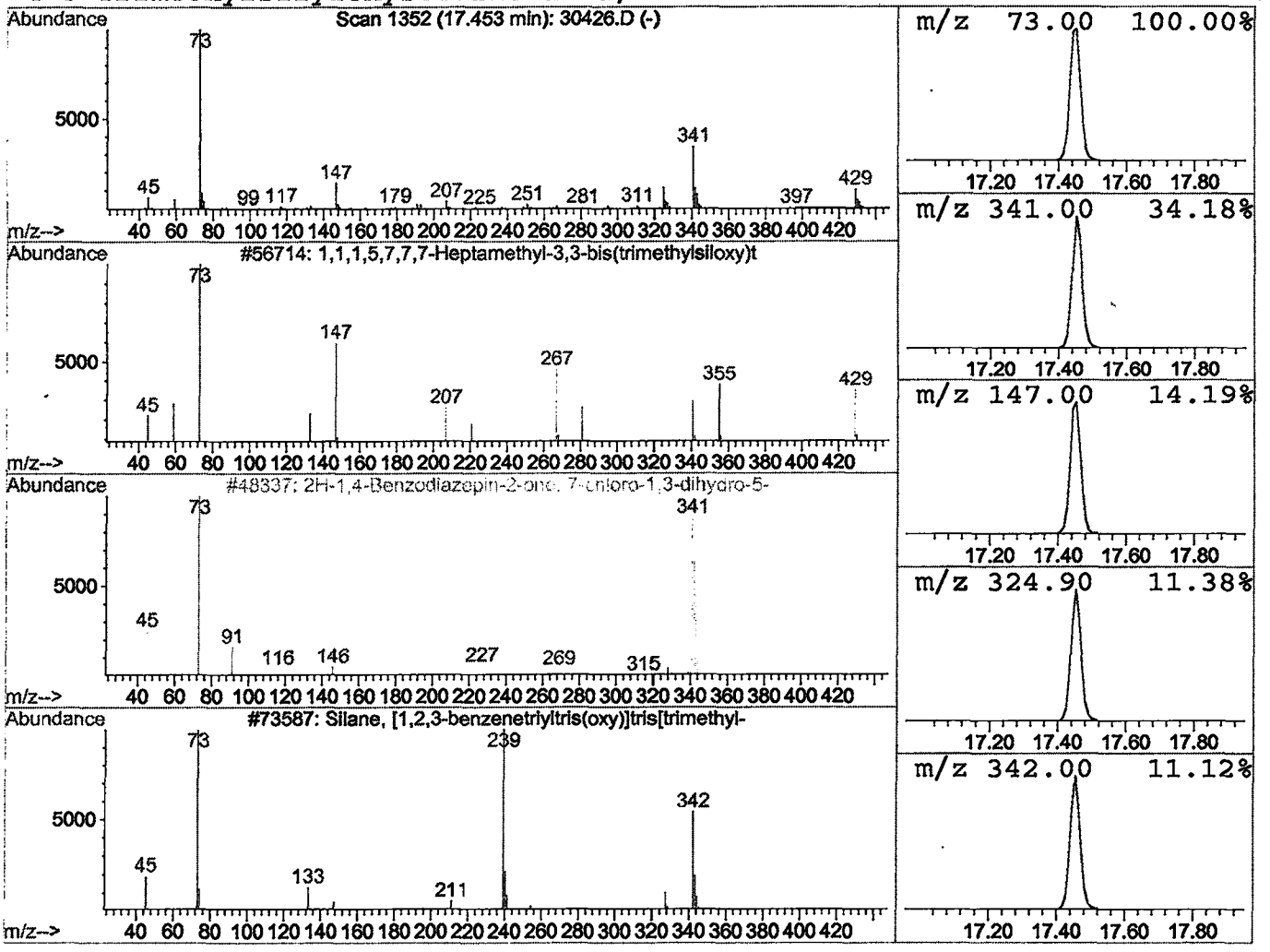
Quant Method : C:\HPCHEM\1\METHODS\GCMS1126.M (RTE Integrator)
Title : 209 625/TCLP calibration table
Library : C:\DATABASE\NBS75K.L

****

Peak Number 4 1,1,1,5,7,7,7-Heptamethyl-3,3- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.45	4.48 ug/l	1870660	Naphthalene-d8	15.28

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	25
2	2H-1,4-Benzodiazepin-2-one, 7-chlor	342	C18H19ClN2OSi	055299-24-6	22
3	Silane, [1,2,3-benzenetriyltris(oxy	342	C15H30O3Si3	017864-23-2	12
4	3-Trimethylsilyloxystearic acid, tr	444	C24H52O3Si2	000000-00-0	9



Library Search Compound Report

Data File : C:\HPCHEM\1\DATA\DEC2701\30426.D
Acq On : 28 Dec 2001 11:28 am
Sample : GC/MS SCAN 12/13
Misc :
MS Integration Params: LSCINT.P

Vial: 23
Operator: 209 GALOS
Inst : GC/MS 209
Multiplr: 1.00

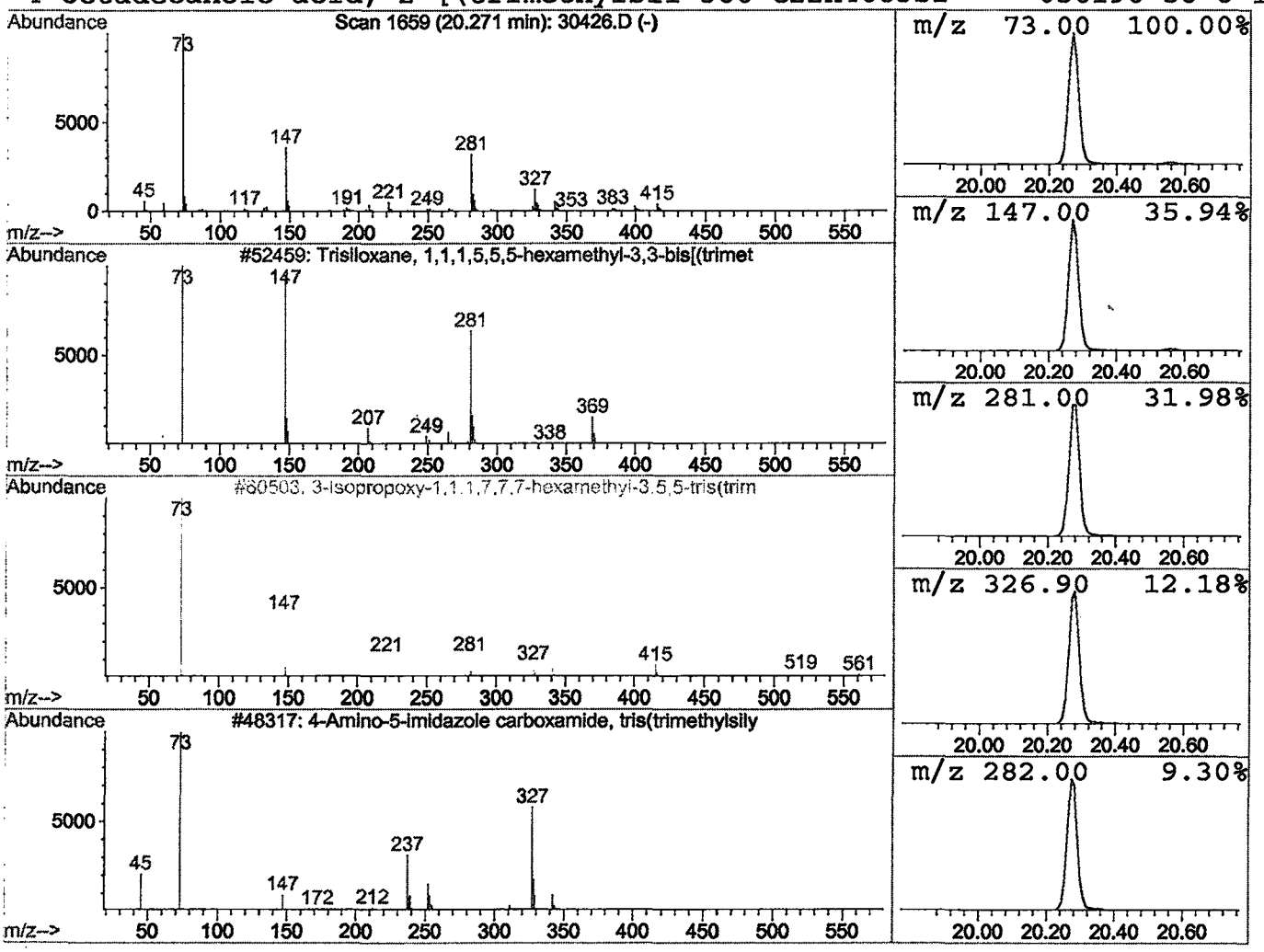
Quant Method : C:\HPCHEM\1\METHODS\GCMS1126.M (RTE Integrator)
Title : 209 625/TCLP calibration table
Library : C:\DATABASE\NBS75K.L

****

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.27	3.47 ug/l	1614360	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	47
2	3-Isopropoxy-1,1,1,7,7,7-hexamethyl	576	C18H52O7Si7	071579-69-6	17
3	4-Amino-5-imidazole carboxamide, tr	342	C13H30N4OSi3	000000-00-0	12
4	Octadecanoic acid, 2-[(trimethylsil	386	C22H46O3Si	056196-58-8	10



Library Search Compound Report

Data File : C:\HPCHEM\1\DATA\DEC2701\30426.D
Acq On : 28 Dec 2001 11:28 am
Sample : GC/MS SCAN 12/13
Misc :
MS Integration Params: LSCINT.P

Vial: 23
Operator: 209 GALOS
Inst : GC/MS 209
Multiplr: 1.00

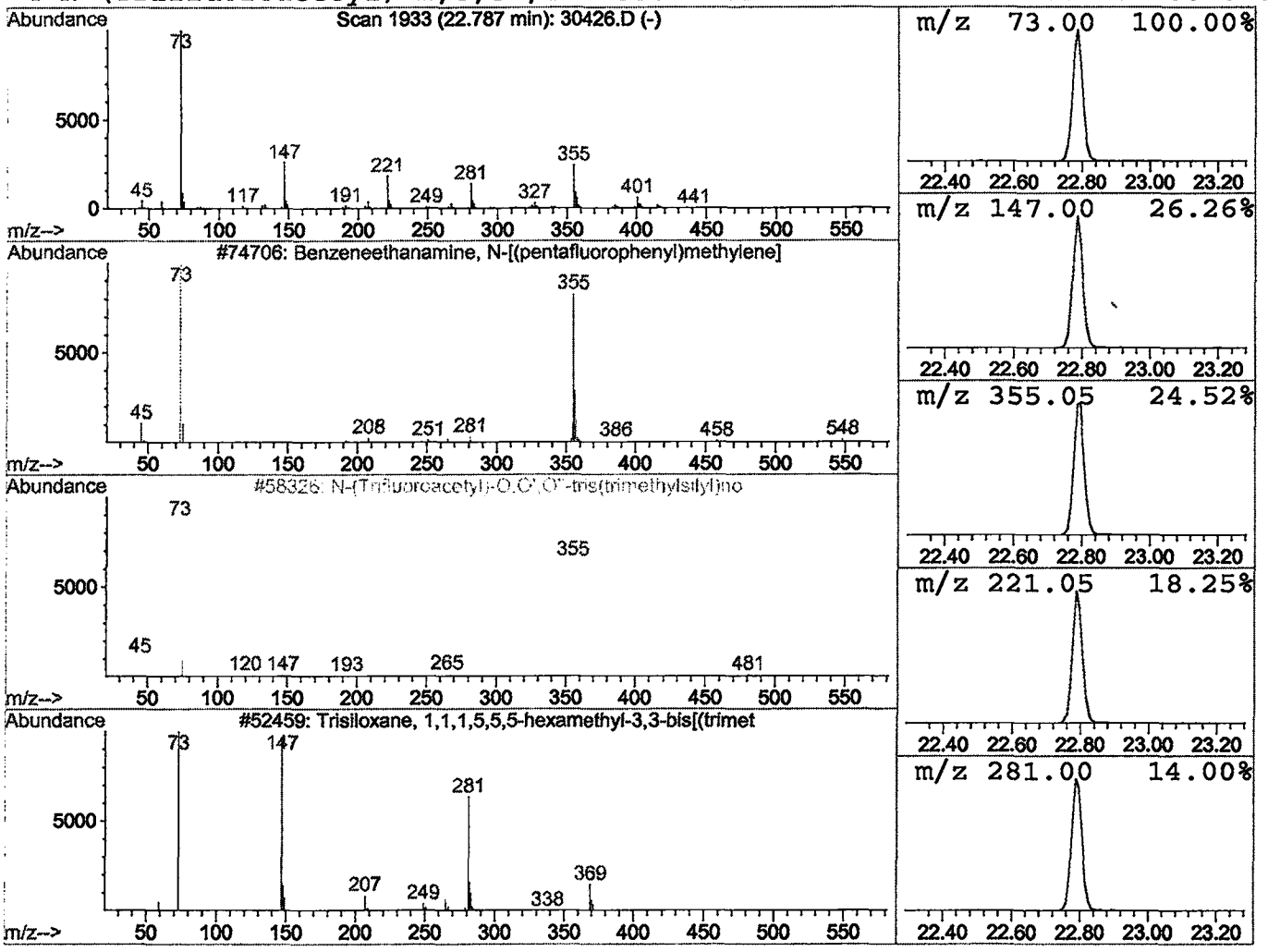
Quant Method : C:\HPCHEM\1\METHODS\GCMS1126.M (RTE Integrator)
Title : 209 625/TCLP calibration table
Library : C:\DATABASE\NBS75K.L

****

Peak Number 6 Benzeneethanamine, N-[(pentafl Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.79	2.70 ug/l	1344980	Phenanthrene-d10	24.98

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzeneethanamine, N-[(pentafluorop	563	C24H34F5NO3Si3	055429-13-5	37
2	N-(Trifluoroacetyl)-O,O',O''-tris(t	481	C19H34F3NO4Si3	000000-00-0	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\DEC2701\30426.D Vial: 23
Acq On : 28 Dec 2001 11:28 am Operator: 209 GALOS
Sample : GC/MS SCAN 12/13 Inst : GC/MS 209
Misc : Multiplr: 1.00
MS Integration Params: LSCINT.P

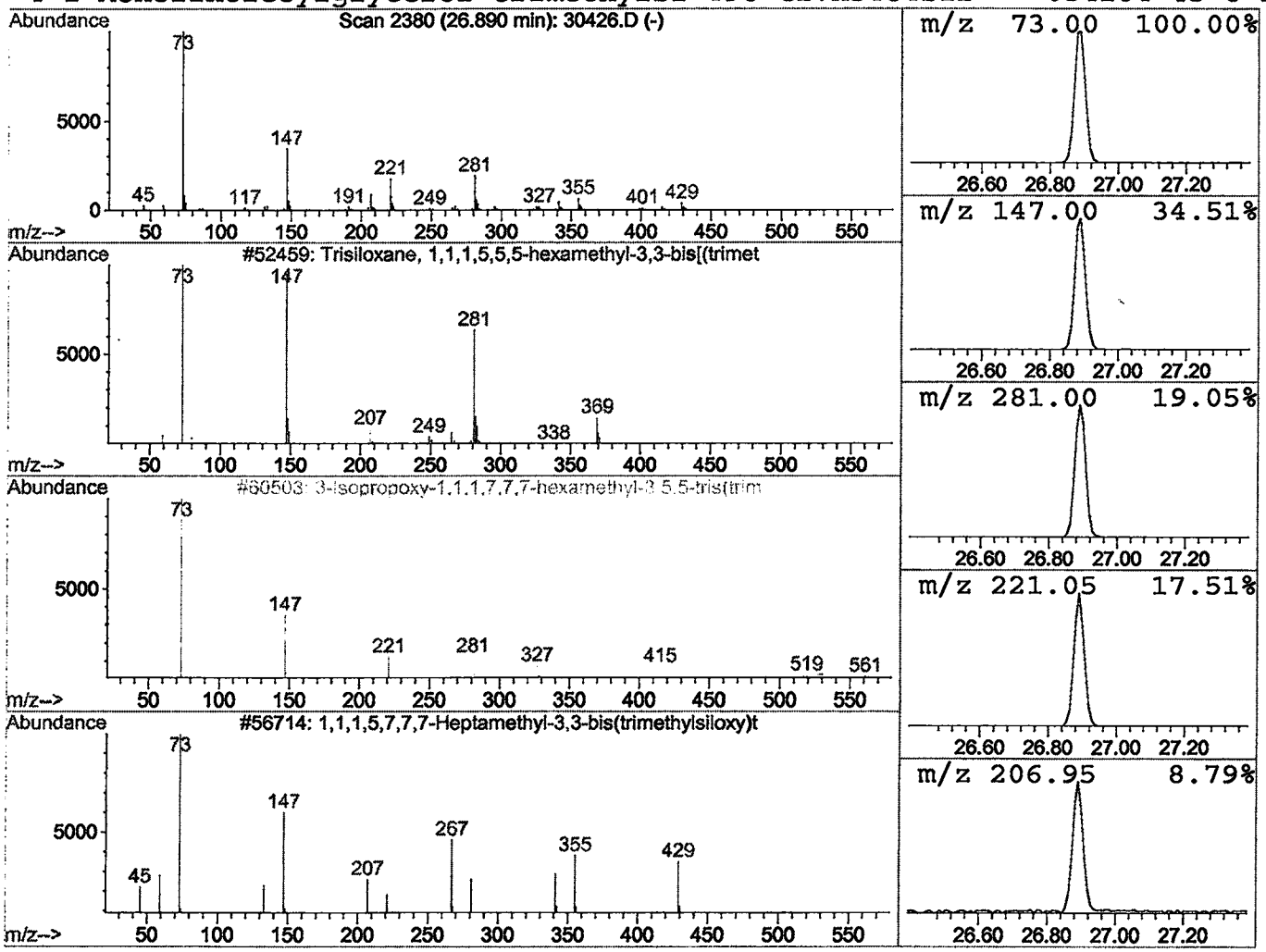
Quant Method : C:\HPCHEM\1\METHODS\GCMS1126.M (RTE Integrator)
Title : 209 625/TCLP calibration table
Library : C:\DATABASE\NBS75K.L

****

Peak Number 7 Trisiloxane, 1,1,1,5,5,5-hexam Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
26.89	1.75 ug/l	869295	Phenanthrene-d10	24.98

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	40
3	1,1,1,5,7,7,7-Heptamethyl-3,3-bis(t	444	C13H40O5Si6	038147-00-1	35
4	1-Monolinoleoylglycerol trimethylsi	498	C27H54O4Si2	054284-45-6	25



Library Search Compound Report

Data File : C:\HPCHEM\1\DATA\DEC2701\30426.D Vial: 23
Acq On : 28 Dec 2001 11:28 am Operator: 209 GALOS
Sample : GC/MS SCAN 12/13 Inst : GC/MS 209
Misc : Multiplr: 1.00
MS Integration Params: LSCINT.P

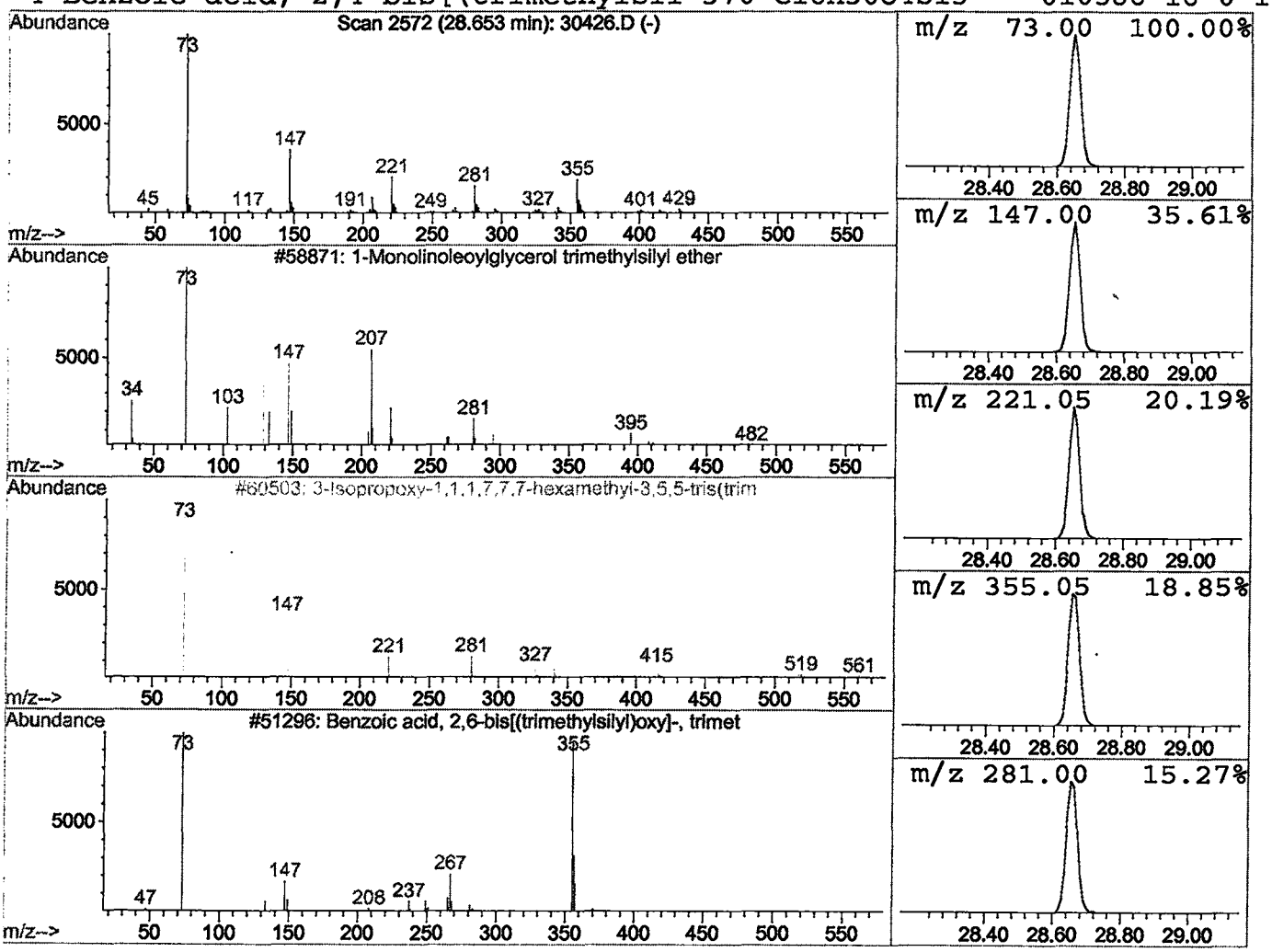
Quant Method : C:\HPCHEM\1\METHODS\GCMS1126.M (RTE Integrator)
Title : 209 625/TCLP calibration table
Library : C:\DATABASE\NBS75K.L

****

Peak Number 8 1-Monolinoleoylglycerol trimet Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
28.65	1.48 ug/l	734867	Phenanthrene-d10	24.98

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Monolinoleoylglycerol trimethylsi	498	C27H54O4Si2	054284-45-6	34
2	3-Isopropoxy-1,1,1,7,7,7-hexamethyl	576	C18H52O7Si7	071579-69-6	25
3	Benzoic acid, 2,6-bis[(trimethylsil	370	C16H30O4Si3	003782-85-2	14
4	Benzoic acid, 2,4-bis[(trimethylsil	370	C16H30O4Si3	010586-16-0	12



Library Search Compound Report

Data File : C:\HPCHEM\1\DATA\DEC2701\30426.D
Acq On : 28 Dec 2001 11:28 am
Sample : GC/MS SCAN 12/13
Misc :
MS Integration Params: LSCINT.P

Vial: 23
Operator: 209 GALOS
Inst : GC/MS 209
Multiplr: 1.00

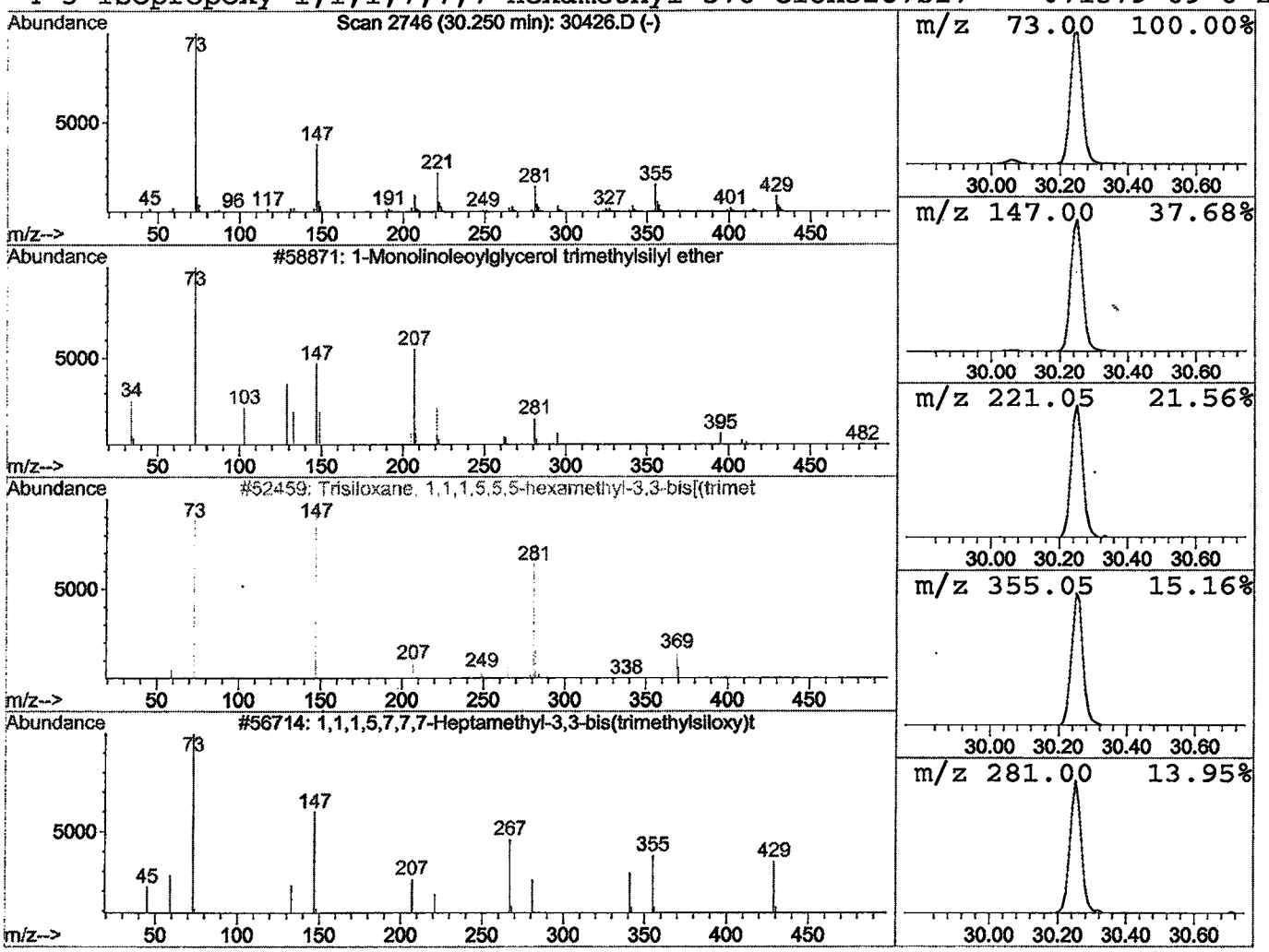
Quant Method : C:\HPCHEM\1\METHODS\GCMS1126.M (RTE Integrator)
Title : 209 625/TCLP calibration table
Library : C:\DATABASE\NBS75K.L

****

Peak Number 9 1-Monolinoleoylglycerol trimet Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
30.25	2.30 ug/l	694473	Chrysene-d12	33.02

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Monolinoleoylglycerol trimethylsi	498	C27H54O4Si2	054284-45-6	42
2	Trisiloxane, 1,1,1,5,5,5-hexamethyl	384	C12H36O4Si5	003555-47-3	32
3	1,1,1,5,7,7,7-Heptamethyl-3,3-bis(t	444	C13H40O5Si6	038147-00-1	28
4	3-Isopropoxy-1,1,1,7,7,7-hexamethyl	576	C18H52O7Si7	071579-69-6	25



Library Search Compound Report

Data File : C:\HPCHEM\1\DATA\DEC2701\30426.D Vial: 23
Acq On : 28 Dec 2001 11:28 am Operator: 209 GALOS
Sample : GC/MS SCAN 12/13 Inst : GC/MS 209
Misc : Multiplr: 1.00
MS Integration Params: LSCINT.P

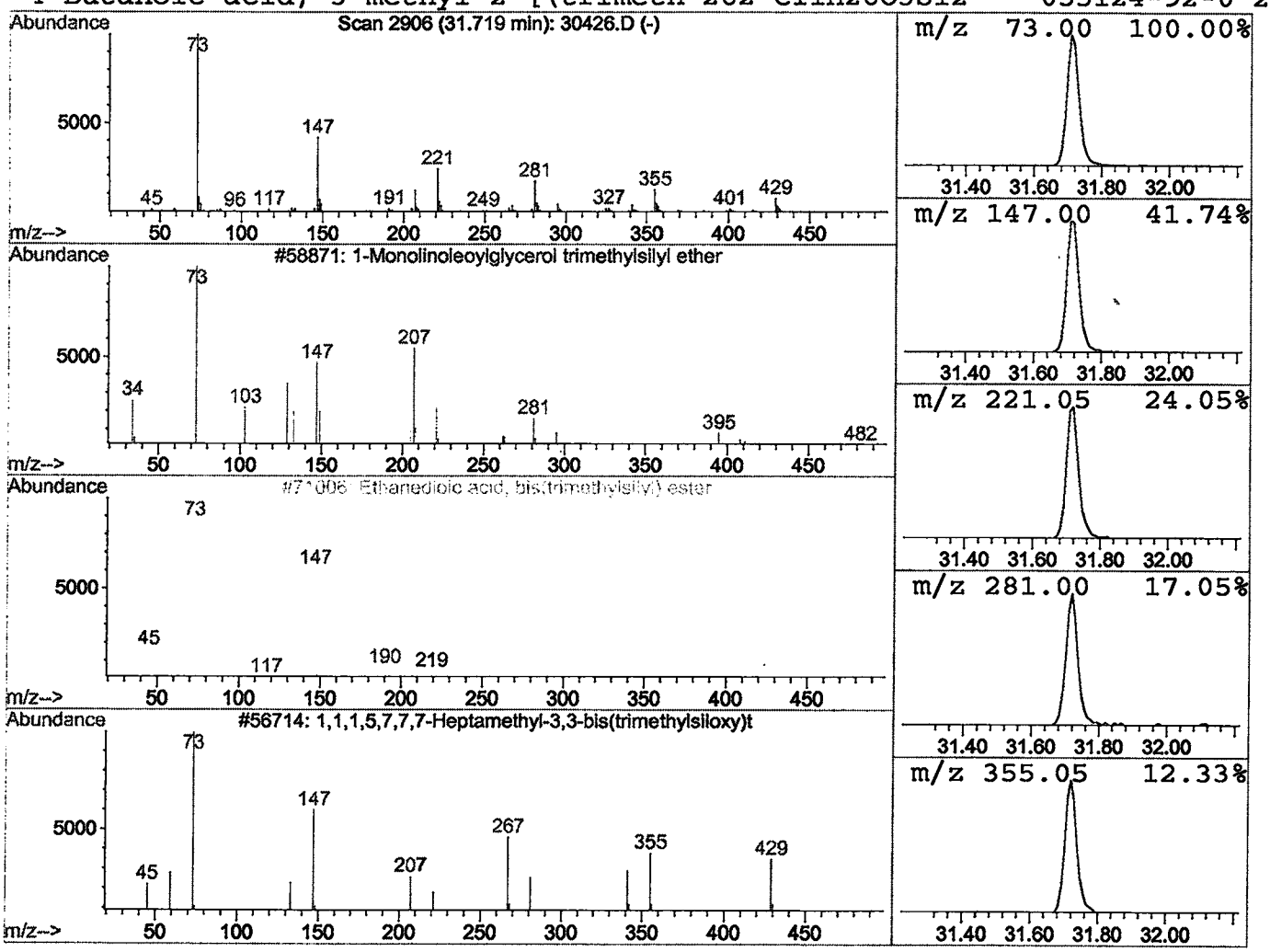
Quant Method : C:\HPCHEM\1\METHODS\GCMS1126.M (RTE Integrator)
Title : 209 625/TCLP calibration table
Library : C:\DATABASE\NBS75K.L

****

Peak Number 10 1-Monolinoleoylglycerol trimet Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
31.72	2.08 ug/l	628540	Chrysene-d12	33.02

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Monolinoleoylglycerol trimethylsi	498	C27H54O4Si2	054284-45-6	45
2	Ethanedioic acid, bis(trimethylsily	234	C8H18O4Si2	018294-04-7	32
3	1,1,1,5,7,7,7-Heptamethyl-3,3-bis(t	444	C13H40O5Si6	038147-00-1	25
4	Butanoic acid, 3-methyl-2-[(trimeth	262	C11H26O3Si2	055124-92-0	23



Library Search Compound Report

Data File : C:\HPCHEM\1\DATA\DEC2701\30426.D
Acq On : 28 Dec 2001 11:28 am
Sample : GC/MS SCAN 12/13
Misc :
MS Integration Params: LSCINT.P

Vial: 23
Operator: 209 GALOS
Inst : GC/MS 209
Multiplr: 1.00

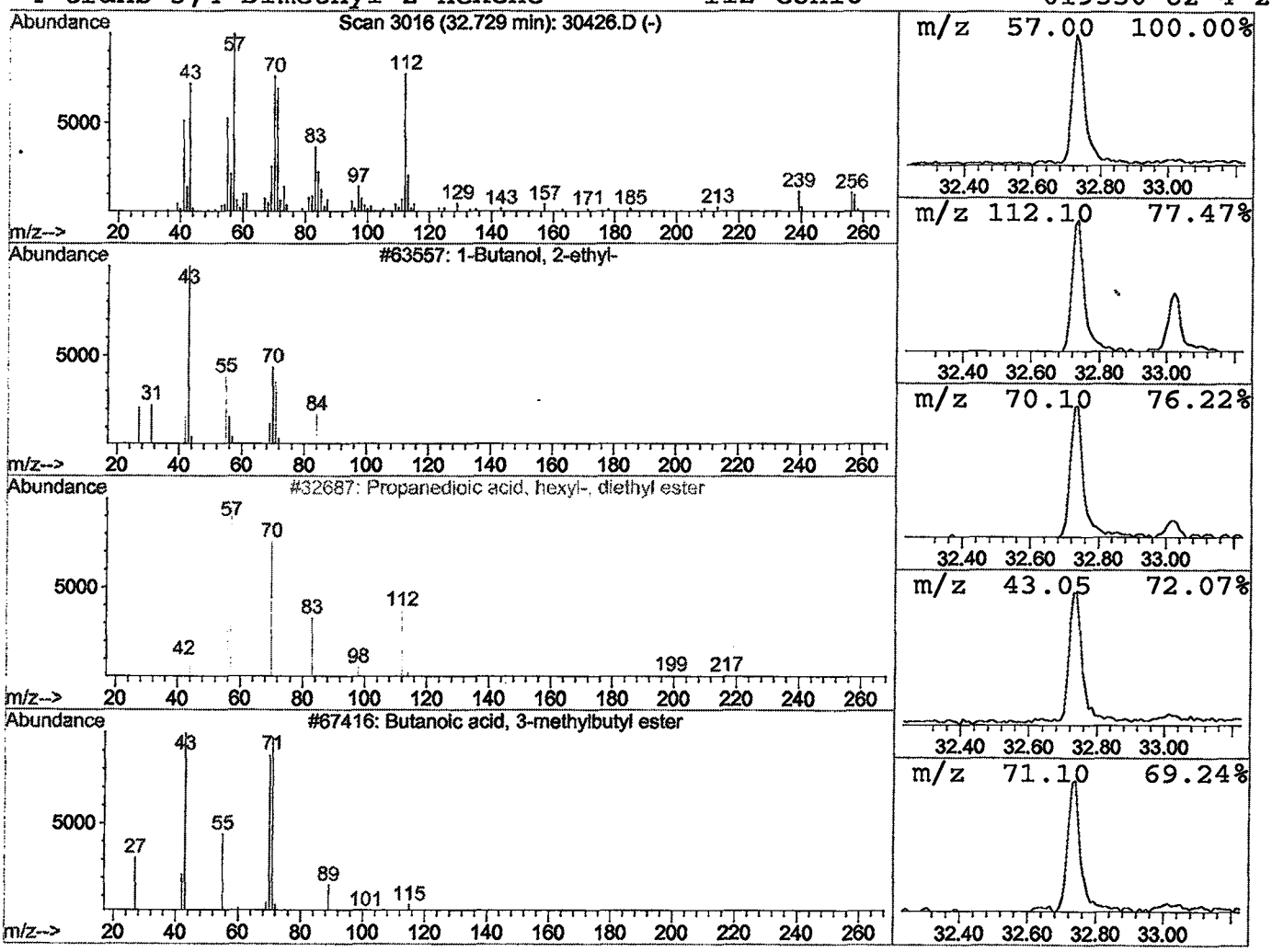
Quant Method : C:\HPCHEM\1\METHODS\GCMS1126.M (RTE Integrator)
Title : 209 625/TCLP calibration table
Library : C:\DATABASE\NBS75K.L

****

Peak Number 11 1-Butanol, 2-ethyl- Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
32.73	0.91 ug/l	274017	Chrysene-d12	33.02

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Butanol, 2-ethyl-	102	C6H14O	000097-95-0	32
2	Propanedioic acid, hexyl-, diethyl	244	C13H24O4	005398-10-7	32
3	Butanoic acid, 3-methylbutyl ester	158	C9H18O2	000106-27-4	27
4	trans-3,4-Dimethyl-2-hexene	112	C8H16	019550-82-4	27



Library Search Compound Report

Data File : C:\HPCHEM\1\DATA\DEC2701\30426.D Vial: 23
Acq On : 28 Dec 2001 11:28 am Operator: 209 GALOS
Sample : GC/MS SCAN 12/13 Inst : GC/MS 209
Misc : Multiplr: 1.00
MS Integration Params: LSCINT.P

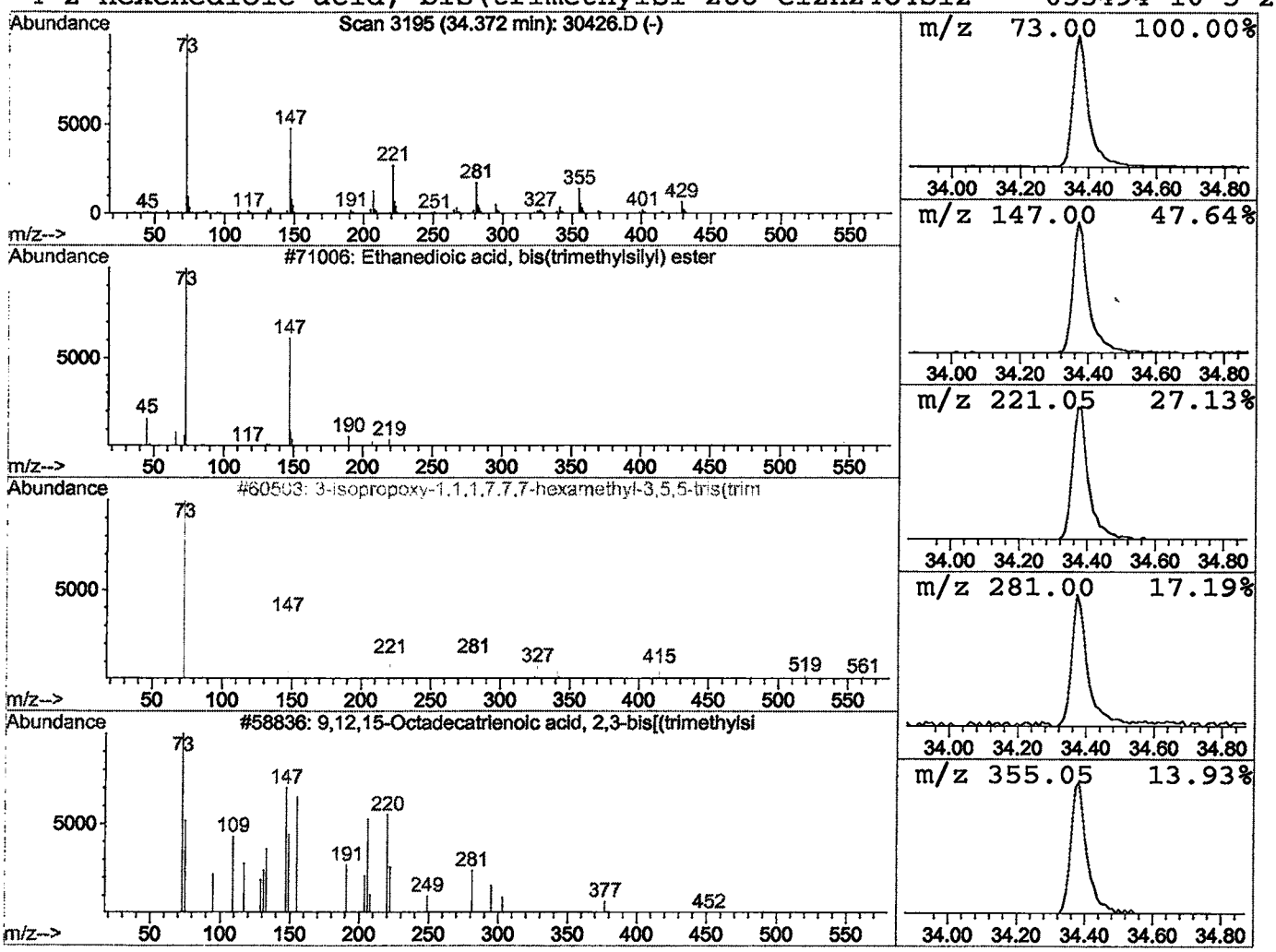
Quant Method : C:\HPCHEM\1\METHODS\GCMS1126.M (RTE Integrator)
Title : 209 625/TCLP calibration table
Library : C:\DATABASE\NBS75K.L

****

Peak Number 12 Ethanedioic acid, bis(trimethy Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
34.37	1.60 ug/l	482977	Chrysene-d12	33.02

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Ethanedioic acid, bis(trimethylsily	234	C8H18O4Si2	018294-04-7	32
2	3-Isopropoxy-1,1,1,7,7,7-hexamethyl	576	C18H52O7Si7	071579-69-6	25
3	9,12,15-Octadecatrienoic acid, 2,3-	496	C27H52O4Si2	055521-22-7	23
4	2-Hexenedioic acid, bis(trimethylsi	288	C12H24O4Si2	055494-10-5	23



Library Search Compound Report

Data File : C:\HPCHEM\1\DATA\DEC2701\30426.D
Acq On : 28 Dec 2001 11:28 am
Sample : GC/MS SCAN 12/13
Misc :
MS Integration Params: LSCINT.P

Vial: 23
Operator: 209 GALOS
Inst : GC/MS 209
Multiplr: 1.00

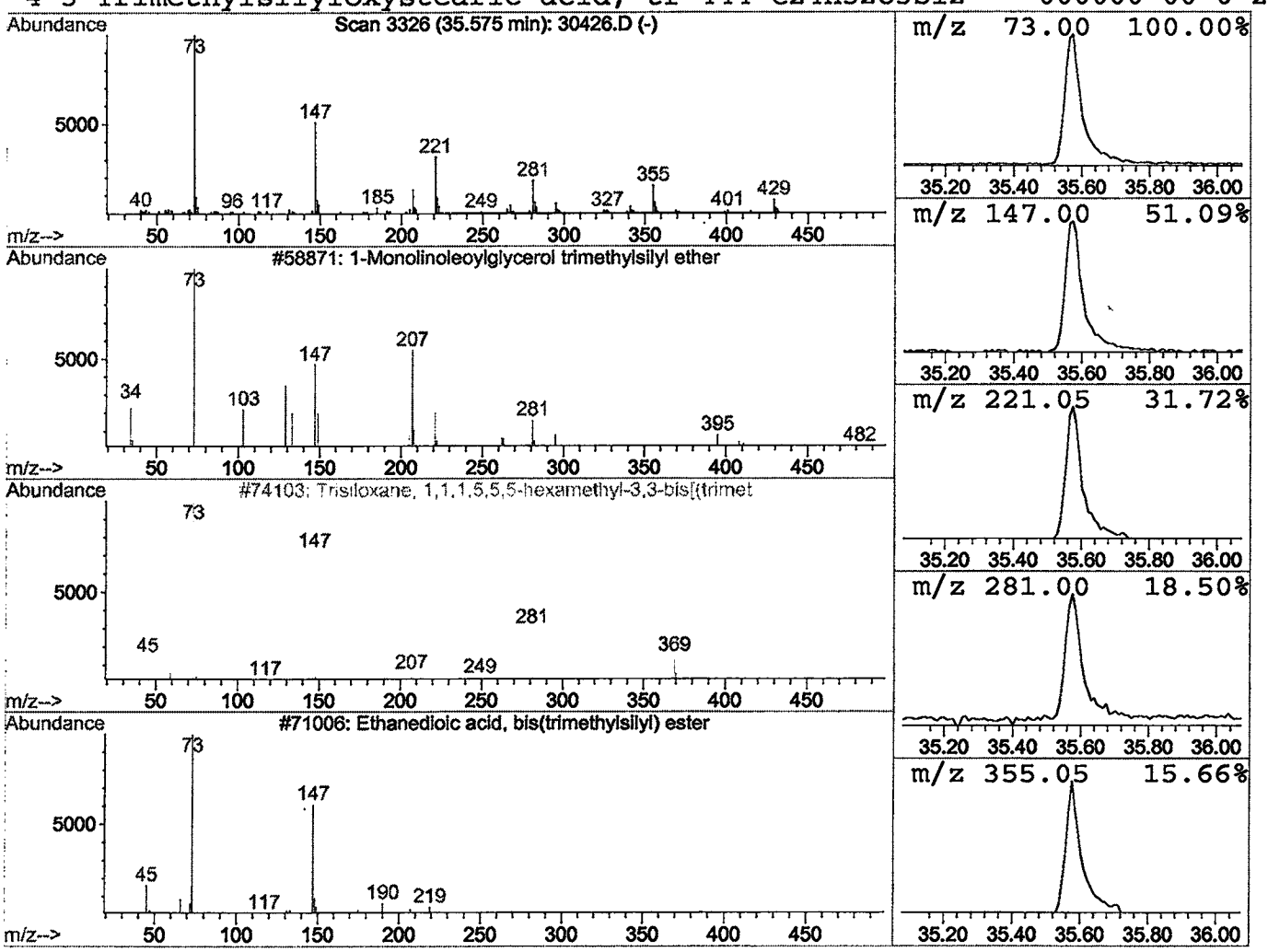
Quant Method : C:\HPCHEM\1\METHODS\GCMS1126.M (RTE Integrator)
Title : 209 625/TCLP calibration table
Library : C:\DATABASE\NBS75K.L

****

Peak Number 13 1-Monolinoleoylglycerol trimet Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
35.57	1.49 ug/l	315476	Perylene-d12	37.12

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Monolinoleoylglycerol trimethylsi	498	C27H54O4Si2	054284-45-6	45
2	Trisiloxane, 1,1,1,5,5,5-hexamethyl	384	C12H36O4Si5	003555-47-3	38
3	Ethanedioic acid, bis(trimethylsily	234	C8H18O4Si2	018294-04-7	38
4	3-Trimethylsilyloxystearic acid, tr	444	C24H52O3Si2	000000-00-0	23



Library Search Compound Report

Data File : C:\HPCHEM\1\DATA\DEC2701\30426.D
Acq On : 28 Dec 2001 11:28 am
Sample : GC/MS SCAN 12/13
Misc :
MS Integration Params: LSCINT.P

Vial: 23
Operator: 209 GALOS
Inst : GC/MS 209
Multiplr: 1.00

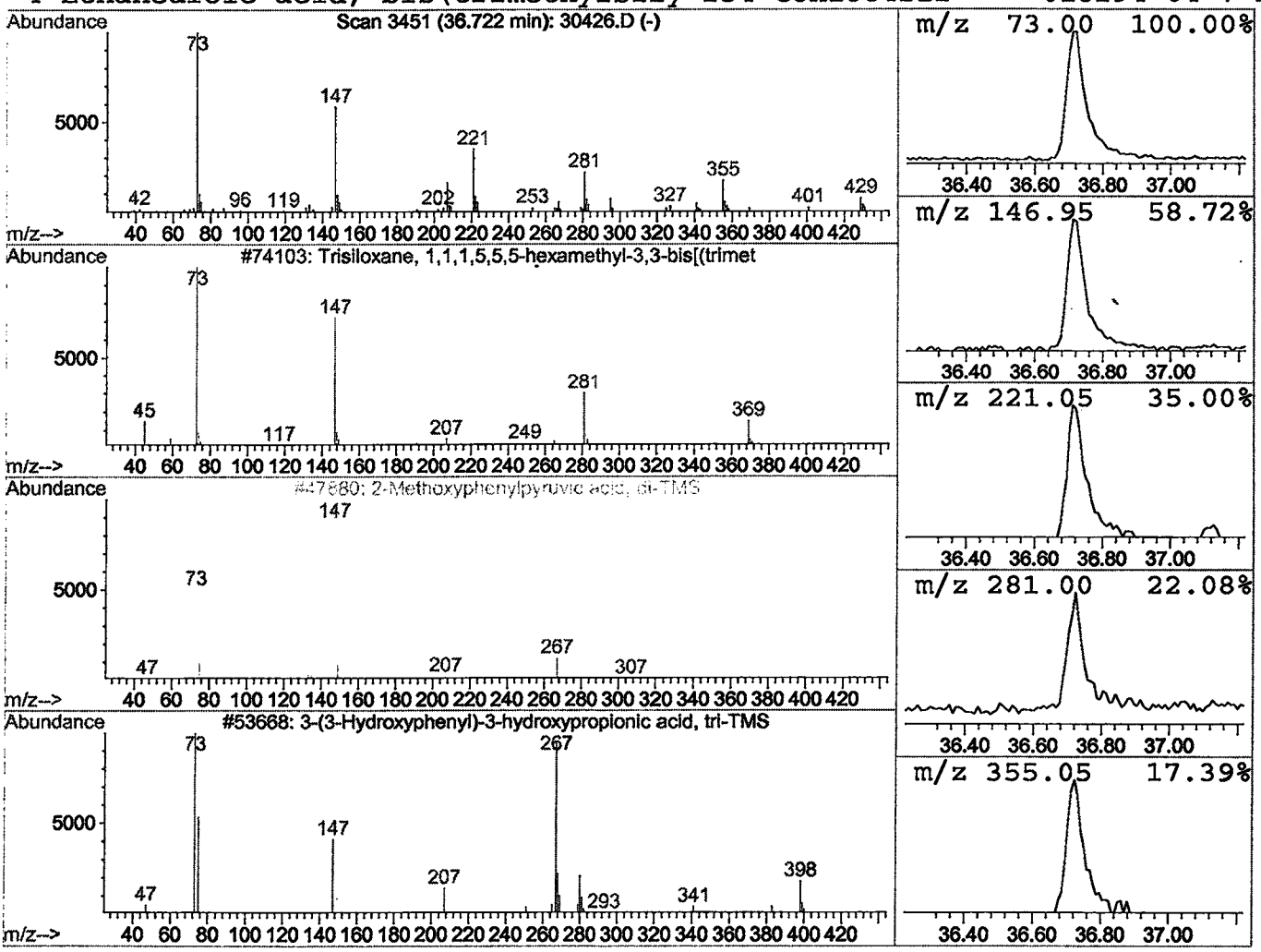
Quant Method : C:\HPCHEM\1\METHODS\GCMS1126.M (RTE Integrator)
Title : 209 625/TCLP calibration table
Library : C:\DATABASE\NBS75K.L

****

Peak Number 14 Trisiloxane, 1,1,1,5,5,5-hexam Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
36.72	0.76 ug/l	161734	Perylene-d12	37.12

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Trisiloxane, 1,1,1,5,5,5-hexamethyl	384	C12H36O4Si5	003555-47-3	38
2	2-Methoxyphenylpyruvic acid, di-TMS	338	C16H26O4Si2	000000-00-0	35
3	3-(3-Hydroxyphenyl)-3-hydroxypropio	398	C18H34O4Si3	000000-00-0	32
4	Ethanedioic acid, bis(trimethylsily	234	C8H18O4Si2	018294-04-7	25



Library Search Compound Report

Data File : C:\HPCHEM\1\DATA\DEC2701\30426.D Vial: 23
Acq On : 28 Dec 2001 11:28 am Operator: 209 GALOS
Sample : GC/MS SCAN 12/13 Inst : GC/MS 209
Misc : Multiplr: 1.00
MS Integration Params: LSCINT.P

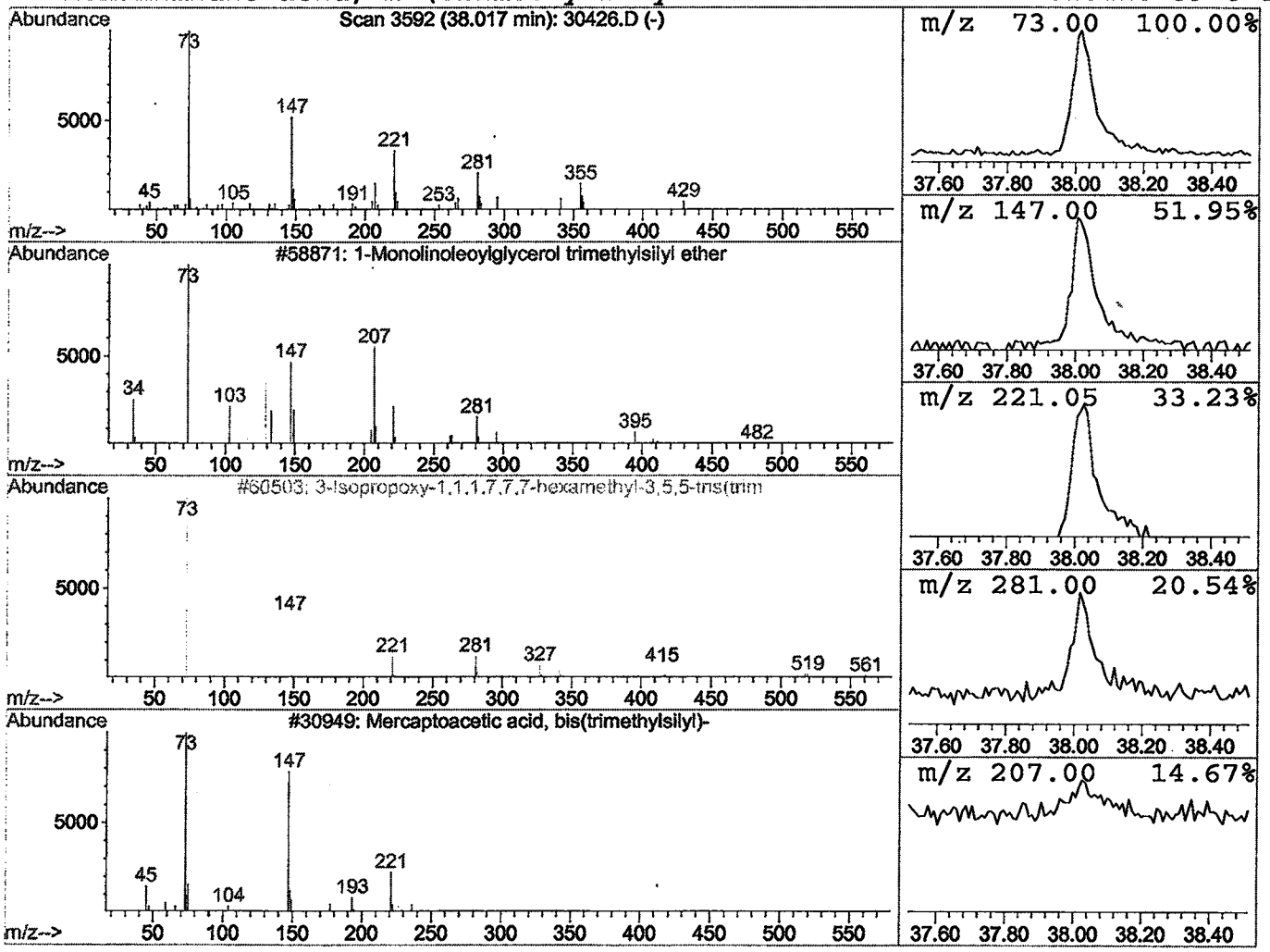
Quant Method : C:\HPCHEM\1\METHODS\GCMS1126.M (RTE Integrator)
Title : 209 625/TCLP calibration table
Library : C:\DATABASE\NBS75K.L

****

Peak Number 15 1-Monolinoleoylglycerol trimet Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
38.02	0.49 ug/l	103388	Perylene-d12	37.12

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Monolinoleoylglycerol trimethylsi	498	C27H54O4Si2	054284-45-6	38
2	3-Isopropoxy-1,1,1,7,7,7-hexamethyl	576	C18H52O7Si7	071579-69-6	28
3	Mercaptoacetic acid, bis(trimethyls	236	C8H20O2SSi2	006398-62-5	17
4	Ethanimidic acid, N-(trimethylsilyl	203	C8H21NOSi2	010416-59-8	12



Tentatively Identified Compound (LSC) summary

Operator ID: 209 GALOS Date Acquired: 28 Dec 2001 11:28 am
Data File: C:\HPCHEM\1\DATA\DEC2701\30426.D
Name: GC/MS SCAN 12/13
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
Silane, tetramethyl-	7.61	4.5	ug/l	1480910	ISTD01	11.68	6639580	20.0
Cyclotetrasiloxane,	11.13	0.5	ug/l	174510	ISTD01	11.68	6639580	20.0
Cyclopentasiloxane,	14.30	2.2	ug/l	929424	ISTD02	15.28	8359550	20.0
1,1,1,5,7,7,7-Heptam	17.45	4.5	ug/l	1870660	ISTD02	15.28	8359550	20.0
Trisiloxane, 1,1,1,5	20.27	3.5	ug/l	1614360	ISTD03	20.56	9309610	20.0
Benzeneethanamine, N	22.79	2.7	ug/l	1344980	ISTD04	24.98	9944590	20.0
Trisiloxane, 1,1,1,5	26.89	1.7	ug/l	869295	ISTD04	24.98	9944590	20.0
1-Monolinoleoylglyce	28.65	1.5	ug/l	734867	ISTD04	24.98	9944590	20.0
1-Monolinoleoylglyce	30.25	2.3	ug/l	694473	ISTD05	33.02	6041610	20.0
1-Monolinoleoylglyce	31.72	2.1	ug/l	628540	ISTD05	33.02	6041610	20.0
1-Butanol, 2-ethyl-	32.73	0.9	ug/l	274017	ISTD05	33.02	6041610	20.0
Ethanedioic acid, bi	34.37	1.6	ug/l	482977	ISTD05	33.02	6041610	20.0
1-Monolinoleoylglyce	35.57	1.5	ug/l	315476	ISTD06	37.12	4231910	20.0
Trisiloxane, 1,1,1,5	36.72	0.8	ug/l	161734	ISTD06	37.12	4231910	20.0
1-Monolinoleoylglyce	38.02	0.5	ug/l	103388	ISTD06	37.12	4231910	20.0

30426.D GCMS1126.M Wed Jan 02 11:08:43 2002