User: Vinci, David L
Westchester Dept. of Labs & Research
Date: 04/26/2002 Time: 14:03:52

Sample: AE09096
Analysis: $GCMS GCMS Scan For Unknowns
Units: ug
Started: 4/26/02 14:02 Ended: 4/26/02 14:02 Analyst: DLV
Page: 1

	Component Name	Vial	Result	Qualifier	MDL
1	Other unknown compounds		.		
2	Surr_2,4-DDD		88		10.0

Comments for Sample AE09096 - Analysis $GCMS

Westchester Dept. of Labs & Research
User: Vinci, David L
Date: 04-26-2002 Time: 14:04:18

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\APR2302\9096.D Vial: 21
Acq On : 24 Apr 2002 10:54 am Operator:
Sample : gc/ms scan 4/17 fb Inst : 209
Misc : Multiplr: 1.00
MS Integration Params: GOOFY.P
Quant Time: Apr 24 12:43 2002 Quant Results File: GCMS0412.RES

Quant Method : C:\HPCHEM\1\METHODS\GCMS0412.M (RTE Integrator)
Title : 209 625/TCLP calibration table
Last Update : Wed Apr 17 11:48:30 2002
Response via : Initial Calibration
DataAcq Meth : GCMS0412

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	12.22	152	468055	20.00	ug/l	-0.02
10) Naphthalene-d8	15.84	136	1729492	20.00	ug/l	-0.03
17) Acenaphthene-d10	21.15	164	971039	20.00	ug/l	-0.02
30) Phenanthrene-d10	25.59	188	1417866	20.00	ug/l	-0.02
39) Chrysene-d12	33.64	240	925683	20.00	ug/l	-0.03
45) Perylene-d12	37.88	264	585695	20.00	ug/l	-0.04
System Monitoring Compounds						
28) TCMX	23.29	209	42680	8.77	ug/L	-0.04
Spiked Amount 10.000			Recovery = 87.70%			
38) 2,4-DDD	0.00	235	0	0.00	ug/mL	
Spiked Amount 5.000			Recovery = 0.00%			
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	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	0.00	128	0	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	0.00	165	0	N.D.		
25) Diethylphthalate	22.63	149	460	N.D.		
26) 4-Chlorophenyl-phenylether	0.00	204	0	N.D.		
27) Fluorene	0.00	166	0	N.D.		
29) Azobenzen/ 1,2-Diphenylhyd	0.00	77	0	N.D.		
31) N-Nitrosodiphenylamine	0.00	168	0	N.D.		
32) 4-Bromophenyl-phenylether	0.00	248	0	N.D.		

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

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Last Update : Wed Apr 17 11:48:30 2002
Response via : Initial Calibration
DataAcq Meth : GCMS0412

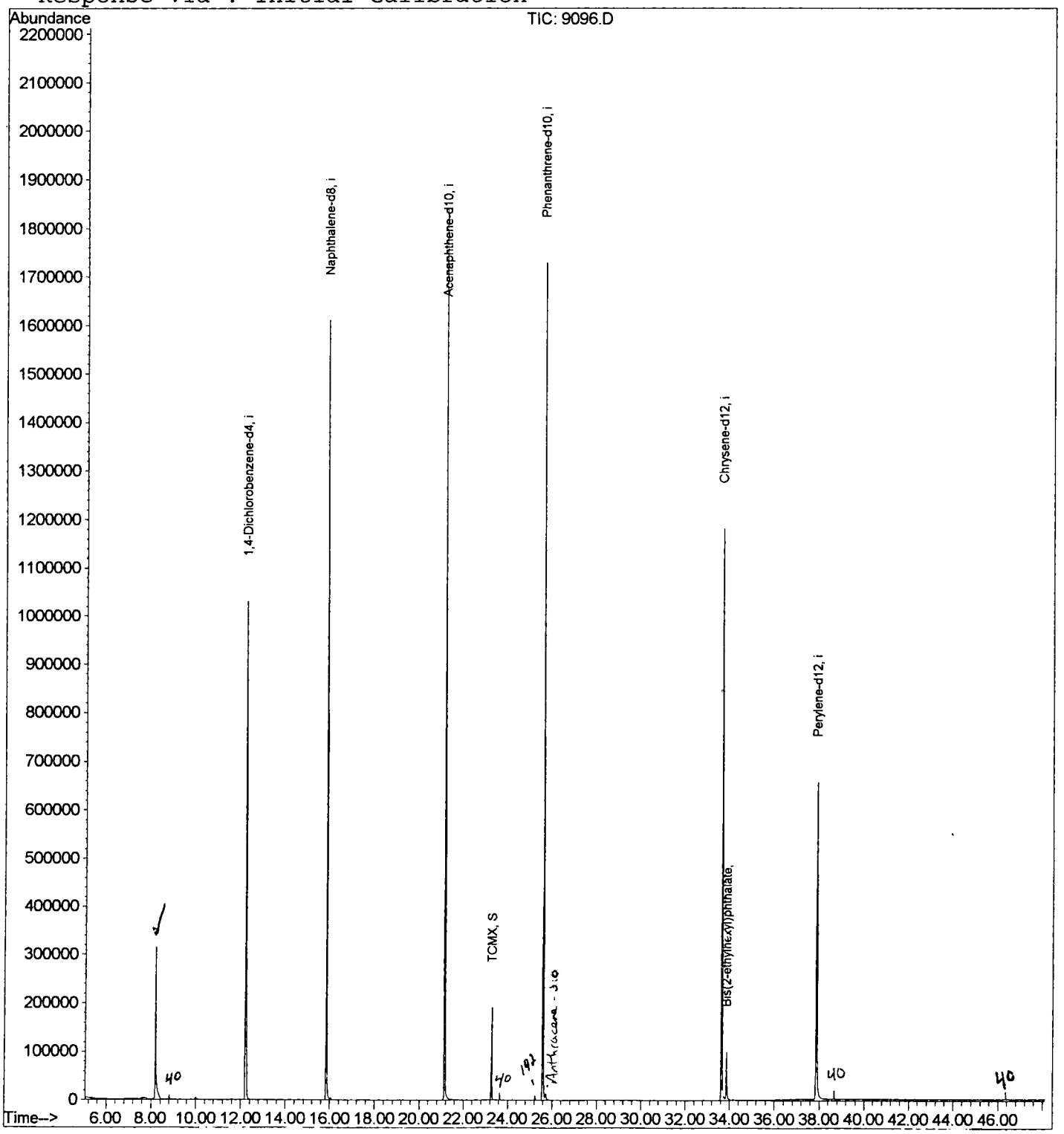
	Compound	R.T.	QIon	Response	Conc Unit	Qvalue
33)	Hexachlorobenzene	0.00	284	0	N.D.	
34)	Phenanthrene	25.58	178	613	N.D.	
35)	Anthracene	25.58	178	612	N.D.	
36)	Di-n-butylphthalate	27.56	149	1497	N.D.	
37)	Fluoranthene	0.00	202	0	N.D.	
40)	Pyrene	0.00	202	0	N.D.	
41)	Butylbenzylphthalate	0.00	149	0	N.D.	
42)	Benzo(a)anthracene	33.65	228	2214	N.D.	
43)	Bis(2-ethylhexyl)phthalate	33.86	149	72921	3.13 ug/l	99
44)	Chrysene	33.65	228	2214	N.D.	
46)	Di-n-Octylphthalate	0.00	149	0	N.D.	
47)	Benzo(b)fluoranthene	0.00	252	0	N.D.	
48)	Benzo(k)fluoranthene	0.00	252	0	N.D.	
49)	Benzo(a)pyrene	37.88	252	2307	N.D.	
50)	Indeno(1,2,3-cd)pyrene	0.00	276	0	N.D.	
51)	Dibenz(a,h)anthracene	0.00	278	0	N.D.	
52)	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\APR2302\9096.D Vial: 21
Acq On : 24 Apr 2002 10:54 am Operator:
Sample : gc/ms scan 4/17 fb Inst : 209
Misc : Multiplr: 1.00
MS Integration Params: GOOFY.P
Quant Time: Apr 24 12:43 2002 Quant Results File: GCMS0412.RES

Method : C:\HPCHEM\1\METHODS\GCMS0412.M (RTE Integrator)
Title : 209 625/TCLP calibration table
Last Update : Wed Apr 17 11:48:30 2002
Response via : Initial Calibration



LSC Area Percent Report

Data File : C:\HPCHEM\1\DATA\APR2302\9096.D Vial: 21
Acq On : 24 Apr 2002 10:54 am Operator:
Sample : gc/ms scan 4/17 fb Inst : 209
Misc : Multiplr: 1.00
MS Integration Params: LSCINT.P

Method : C:\HPCHEM\1\METHODS\GCMS0412.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

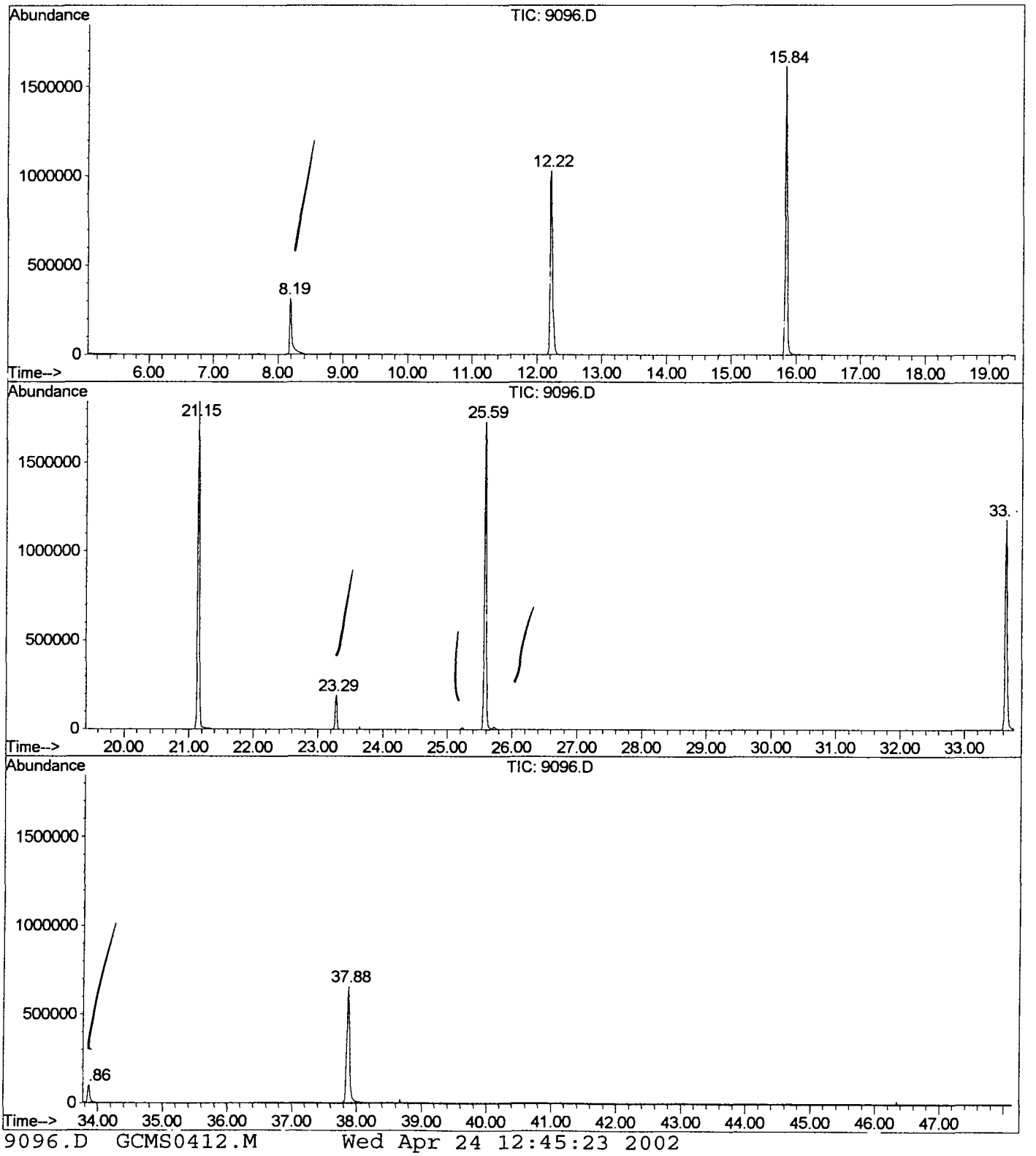
peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	peak area	peak % max.	% of total
1	8.187	338	342	371	rVB	313563	762742	20.50%	3.970%
2	12.217	775	781	802	rBV	1031032	2542946	68.35%	13.235%
3	15.843	1170	1176	1194	rBV	1611241	3278654	88.13%	17.065%
4	21.149	1748	1754	1768	rBV	1843515	3720220	100.00%	19.363%
5	23.288	1979	1987	1994	rVB	192032	380232	10.22%	1.979%
6	25.593	2231	2238	2249	rBV	1730664	3620059	97.31%	18.842%
7	33.644	3108	3115	3134	rBV2	1183786	2693989	72.41%	14.022%
8	33.864	3134	3139	3149	rVV	98159	228816	6.15%	1.191%
9	37.876	3568	3576	3594	rBV2	655117	1985531	53.37%	10.334%

Sum of corrected areas: 19213189

9096.D GCMS0412.M Wed Apr 24 12:45:23 2002

LSC Report - Integrated Chromatogram

File : C:\HPCHEM\1\DATA\APR2302\9096.D
Operator :
Acquired : 24 Apr 2002 10:54 am using AcqMethod GCMS0412
Instrument : 209
Sample Name: gc/ms scan 4/17 fb
Misc Info :
Vial Number: 21
Quant File :GCMS0412.RES (RTE Integrator)



9096.D GCMS0412.M Wed Apr 24 12:45:23 2002

Library Search Compound Report

Data File : C:\HPCHEM\1\DATA\APR2302\9096.D Vial: 21
Acq On : 24 Apr 2002 10:54 am Operator:
Sample : gc/ms scan 4/17 fb Inst : 209
Misc : Multiplr: 1.00
MS Integration Params: LSCINT.P

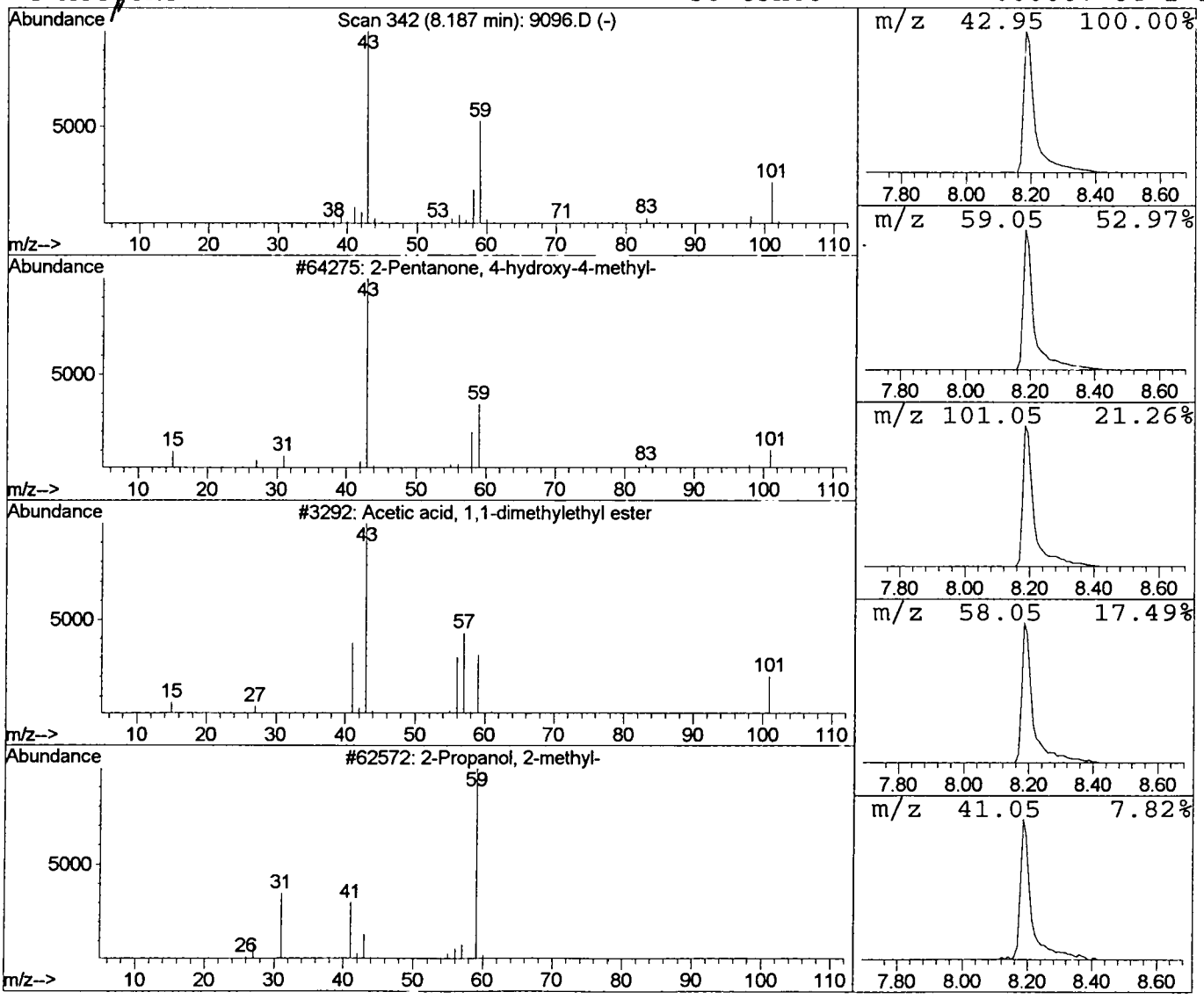
Quant Method : C:\HPCHEM\1\METHODS\GCMS0412.M (RTE Integrator)
Title : 209 625/TCLP calibration table
Library : C:\DATABASE\NBS75K.L

****

Peak Number 1 2-Pentanone, 4-hydroxy-4-methy Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.19	6.00 ug/l	762742	1,4-Dichlorobenzene-d4	12.22

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	50
2	Acetic acid, 1,1-dimethylethyl este	116	C6H12O2	000540-88-5	33
3	2-Propanol, 2-methyl-	74	C4H10O	000075-65-0	17
4	Acetone	58	C3H6O	000067-64-1	10



Tentatively Identified Compound (LSC) summary

Operator ID: Date Acquired: 24 Apr 2002 10:54 am
Data File: C:\HPCHEM\1\DATA\APR2302\9096.D
Name: gc/ms scan 4/17 fb
Misc:
Method: C:\HPCHEM\1\METHODS\GCMS0412.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
~~2-Pentanone, 4-hydro~~	8.19	6.0	ug/l	762742	ISTD01	12.22	2542950	20.0

9096.D GCMS0412.M Wed Apr 24 12:45:25 2002