User: Vinci, David L
Westchester Dept. of Labs & Research
Date: 04/29/2002 Time: 15:01:04

Sample: AE07068
Analysis: $GCMS GCMS Scan For Unknowns
Units: ug
Started: 4/29/02 15:00 Ended: 4/29/02 15:00 Analyst: DLV
Page: 1

	Component Name	Viol	Result	Qualifier	MDL
1	Other unknown compounds		80		
2	Surr_2,4-DDD				10.0

User: Vinci, David L
Westchester Dept. of Labs & Research
Date: 04/29/2002 Time: 15:01:14

Sample: AE07068
Analysis: $GCMS GCMS Scan For Unknowns
Units: ug
Started: 4/29/02 15:00 Ended: 4/29/02 15:00 Analyst: DLV
Page: 1

	Component Name	Viol	Result	Qualifier	MDL
1	Other unknown compounds		.		
2	Surr_2,4-DDD		80		10.0

Comments for Sample AE07068 - Analysis $GCMS

Westchester Dept. of Labs & Research
User: Vinci, David L
Date: 04-29-2002 Time: 15:01:39

The GCMS scan, from 35 to 550 amu, does not reveal the presence of unusual compounds.
The recoveries for 5 of the 43 compounds in the LCS Standard were low.

Quantitation Report (QT Reviewed)

Data File : C:\HPCHEM\1\DATA\APR2402\7068.D Vial: 10
Acq On : 25 Apr 2002 12:34 am Operator:
Sample : gc/ms scan 4/23 fb Inst : 209
Misc : Multiplr: 1.00
MS Integration Params: GOOFY.P
Quant Time: Apr 25 8:31 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	12.21	152	530365	20.00	ug/l	-0.03
10) Naphthalene-d8	15.84	136	1929775	20.00	ug/l	-0.03
17) Acenaphthene-d10	21.15	164	1082802	20.00	ug/l	-0.02
30) Phenanthrene-d10	25.58	188	1567770	20.00	ug/l	-0.03
39) Chrysene-d12	33.64	240	1045943	20.00	ug/l	-0.03
45) Perylene-d12	37.87	264	682329	20.00	ug/l	-0.05
System Monitoring Compounds						
28) TCMX	23.28	209	43365	7.99	ug/L	-0.05
Spiked Amount 10.000			Recovery = 79.90%			
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	15.90	128	185	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	22.62	149	836	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

Quantitation Report (QT Reviewed)

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Sample : gc/ms scan 4/23 fb Inst : 209
Misc : Multiplr: 1.00
MS Integration Params: GOOFY.P
Quant Time: Apr 25 8:31 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

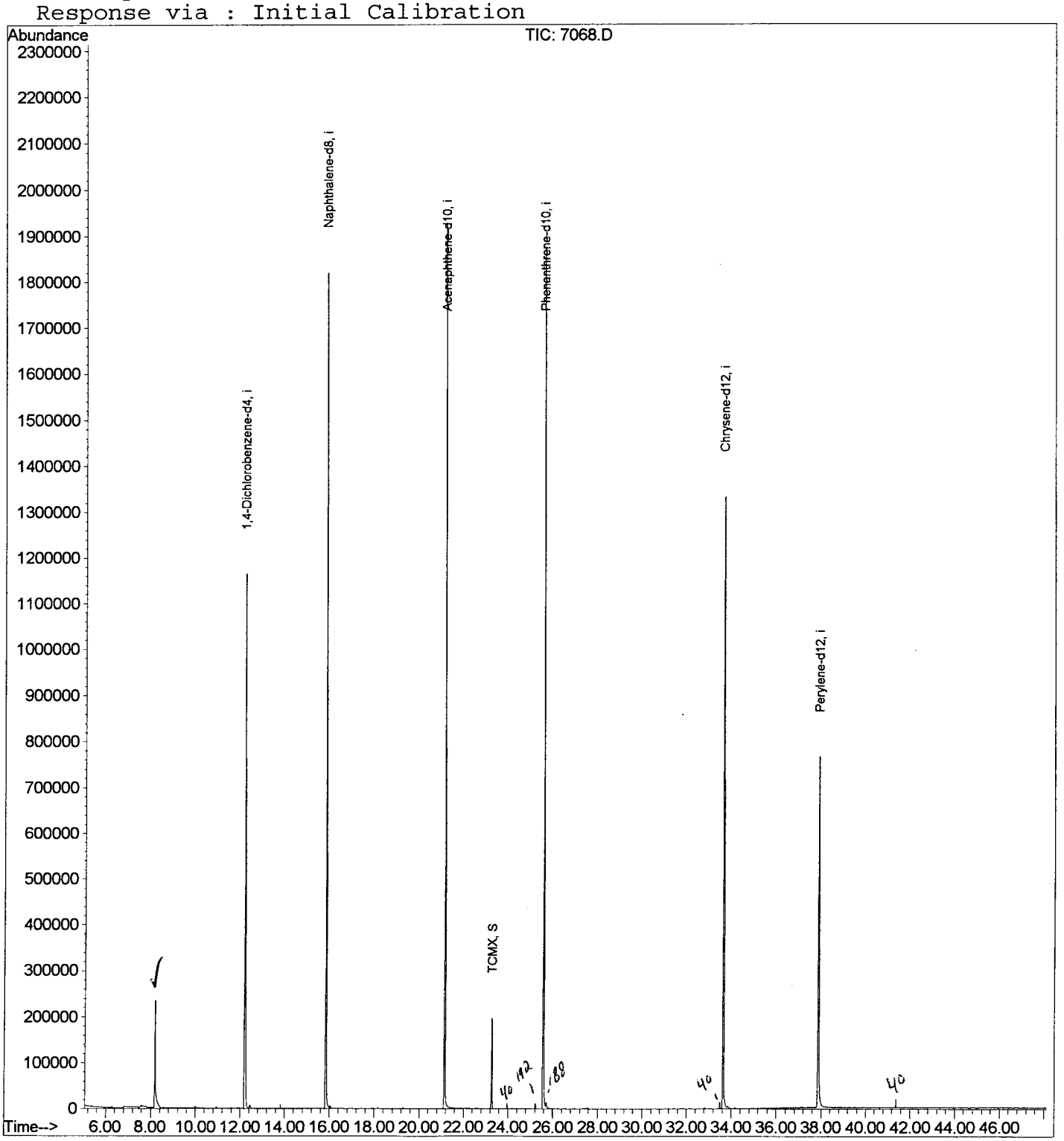
	Compound	R.T.	QIon	Response	Conc Unit	Qvalue
33)	Hexachlorobenzene	0.00	284	0	N.D.	
34)	Phenanthrene	25.59	178	485	N.D.	
35)	Anthracene	25.59	178	485	N.D.	
36)	Di-n-butylphthalate	27.56	149	4144	N.D.	
37)	Fluoranthene	0.00	202	0	N.D.	
40)	Pyrene	0.00	202	0	N.D.	
41)	Butylbenzylphthalate	0.00	149	0	N.D.	
42)	Benzo(a)anthracene	33.64	228	2262	N.D.	
43)	Bis(2-ethylhexyl)phthalate	33.86	149	2317	N.D.	
44)	Chrysene	33.64	228	2262	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	2736	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\APR2402\7068.D Vial: 10
Acq On : 25 Apr 2002 12:34 am Operator:
Sample : gc/ms scan 4/23 fb Inst : 209
Misc : Multiplr: 1.00
MS Integration Params: GOOFY.P
Quant Time: Apr 25 8:31 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\APR2402\7068.D Vial: 10
Acq On : 25 Apr 2002 12:34 am Operator:
Sample : gc/ms scan 4/23 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

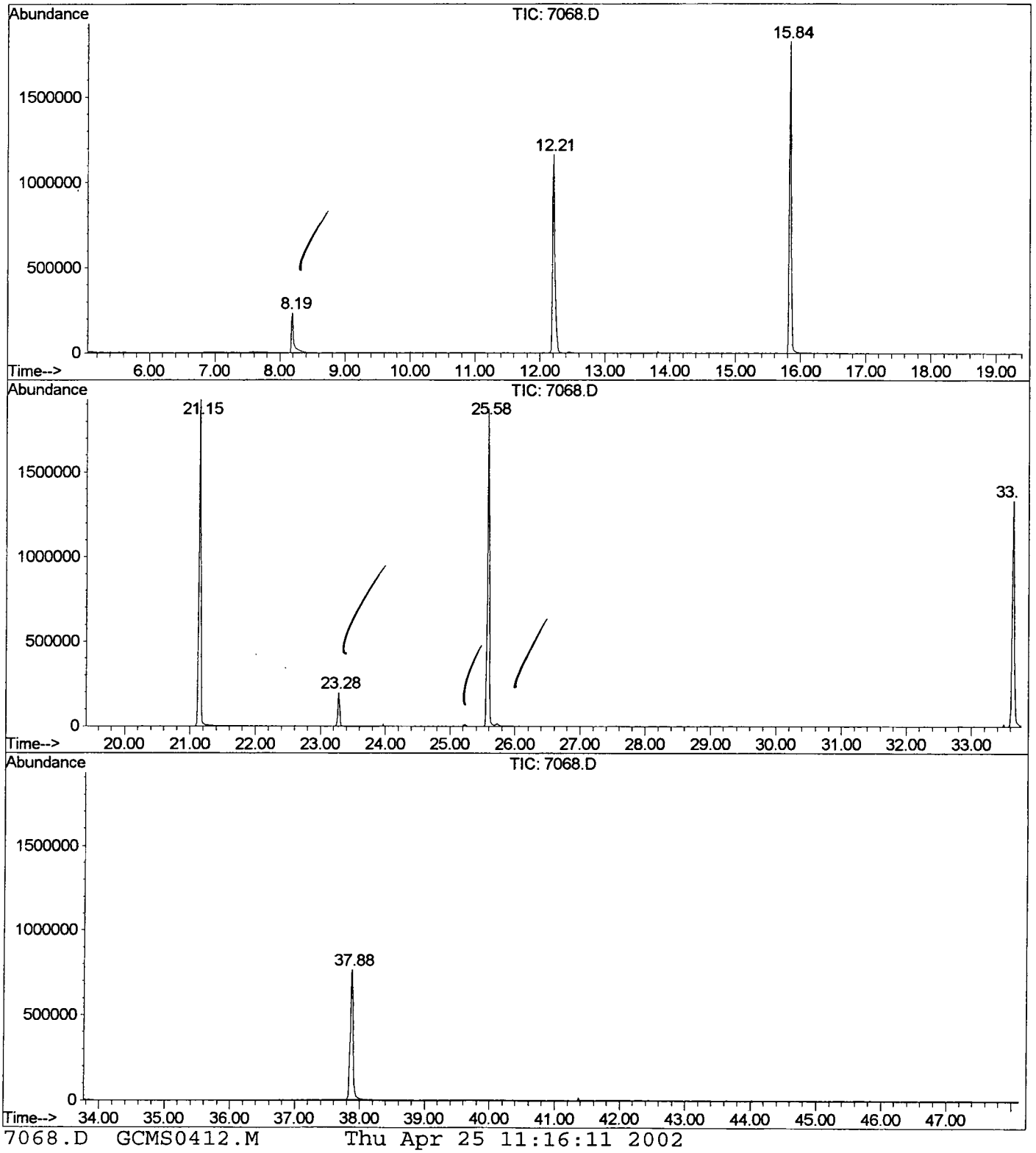
peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	peak area	peak % max.	% of total
1	8.187	338	342	371	rVB	233146	577898	14.15%	2.779%
2	12.208	775	780	798	rVB	1163151	2853160	69.85%	13.720%
3	15.844	1170	1176	1190	rBV	1820240	3647766	89.31%	17.541%
4	21.150	1747	1754	1765	rBV	1928298	4084615	100.00%	19.642%
5	23.280	1978	1986	1993	rBV	197900	387386	9.48%	1.863%
6	25.584	2230	2237	2248	rBV2	1872918	3927732	96.16%	18.887%
7	33.644	3108	3115	3132	rBV	1332661	3007305	73.63%	14.461%
8	37.877	3567	3576	3600	rBV2	763763	2309755	56.55%	11.107%

Sum of corrected areas: 20795617

7068.D GCMS0412.M Thu Apr 25 11:16:10 2002

LSC Report - Integrated Chromatogram

File : C:\HPCHEM\1\DATA\APR2402\7068.D
Operator :
Acquired : 25 Apr 2002 12:34 am using AcqMethod GCMS0412
Instrument : 209
Sample Name: gc/ms scan 4/23 fb
Misc Info :
Vial Number: 10
Quant File :GCMS0412.RES (RTE Integrator)



7068.D GCMS0412.M Thu Apr 25 11:16:11 2002

Library Search Compound Report

Data File : C:\HPCHEM\1\DATA\APR2402\7068.D Vial: 10
Acq On : 25 Apr 2002 12:34 am Operator:
Sample : gc/ms scan 4/23 fb Inst : 209
Misc : Multiplr: 1.00
MS Integration Params: LSCINT.P

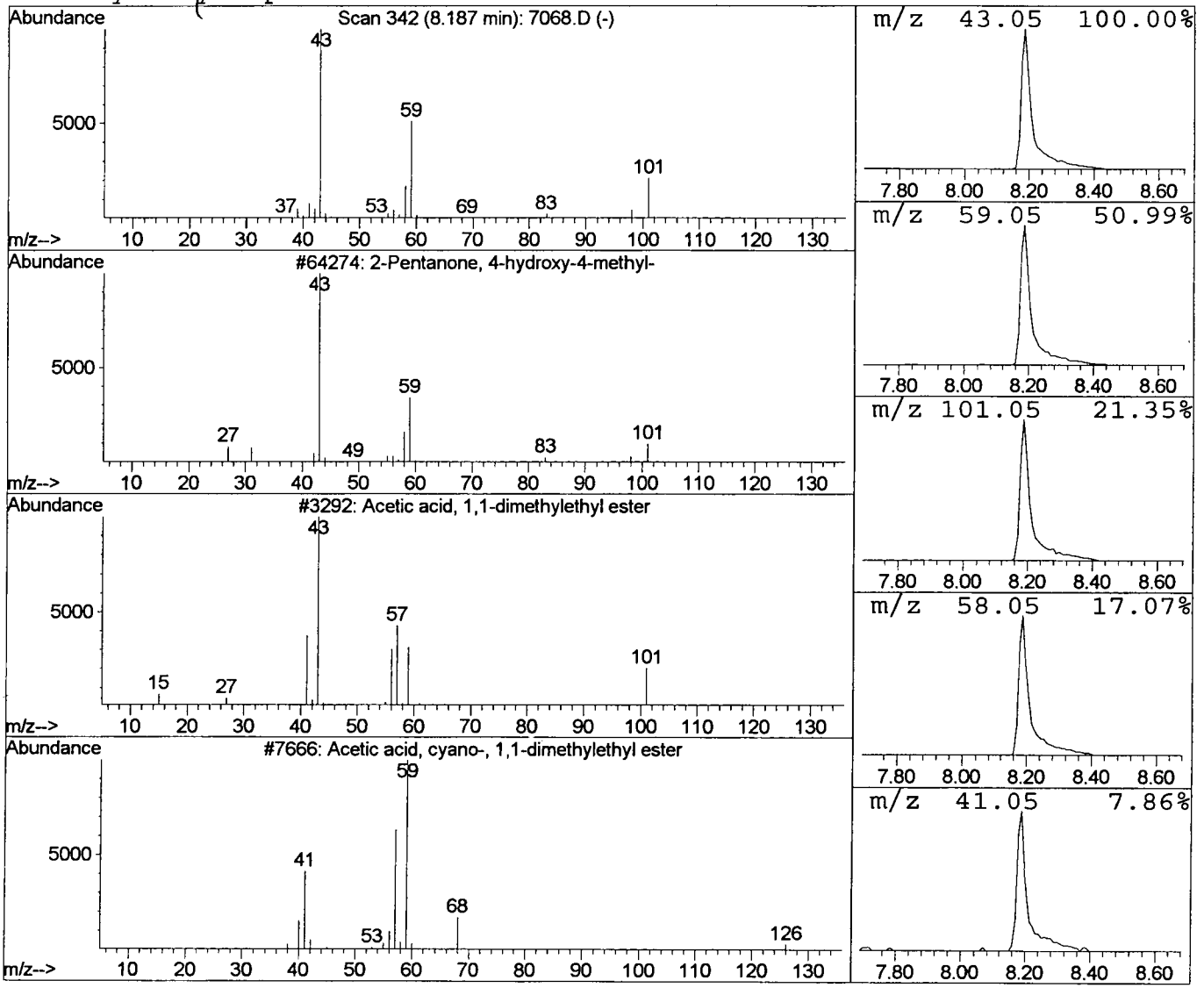
Quant Method : C:\HPCHEM\1\METHODS\GCMS0412.M (RTE Integrator)
Title : 209 625/TCLP calibration table
Library : C:\DATABASE\NBS75K.L

****

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.19	4.05 ug/l	577898	1,4-Dichlorobenzene-d4	12.21

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	72
2	Acetic acid, 1,1-dimethylethyl este	116	C6H12O2	000540-88-5	33
3	Acetic acid, cyano-, 1,1-dimethylet	141	C7H11NO2	001116-98-9	17
4	3-Hydroxy-2-pentanone	102	C5H10O2	003142-66-3	9



Tentatively Identified Compound (LSC) summary

Operator ID: Date Acquired: 25 Apr 2002 12:34 am
Data File: C:\HPCHEM\1\DATA\APR2402\7068.D
Name: gc/ms scan 4/23 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	4.1	ug/l	577898	ISTD01	12.21	2853160	20.0

7068.D GCMS0412.M Thu Apr 25 11:16:12 2002