User: Vinci, David L
Westchester Dept. of Labs & Research
Date: 06/06/2002 Time: 15:39:36

Sample: AE12154
Analysis: $GCMS GCMS Scan For Unknowns
Units: ug
Started: 6/6/02 15:39 Ended: 6/6/02 15:39 Analyst: DLV
Page: 1

	Component Name	Viol	Result	Qualifier	MDL
1	Other unknown compounds		.		
2	Surr_2,4-DDD		86		10.0

Comments for Sample AE12154 - Analysis $GCMS

Westchester Dept. of Labs & Research
User: Vinci, David L
Date: 06-06-2002 Time: 15:42:36

The GCMS scan, from 35 to 550 amu, does not reveal the presence of unusual compounds. The recovery for 1 of the 43 compounds in the LCS Standard was low.

Quantitation Report (QT Reviewed)

Data File : C:\HPCHEM\1\DATA\MAY2402\12154.D Vial: 8
Acq On : 24 May 2002 10:26 pm Operator: Morgan
Sample : gc/ms scan 5/22 fb Inst : 209
Misc : Multiplr: 1.00
MS Integration Params: GOOFY.P
Quant Time: May 28 11:47 2002 Quant Results File: GCMS0520.RES

Quant Method : C:\HPCHEM\1\METHODS\GCMS0520.M (RTE Integrator)
Title : 209 625/TCLP calibration table
Last Update : Tue May 28 11:39:13 2002
Response via : Initial Calibration
DataAcq Meth : GCMS0520

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	12.14	152	527736	40.00	ug/l	0.00
10) Naphthalene-d8	15.78	136	2115431	40.00	ug/l	-0.01
17) Acenaphthene-d10	21.09	164	1010203	40.00	ug/l	0.00
30) Phenanthrene-d10	25.52	188	1439676	40.00	ug/l	-0.01
38) Chrysene-d12	33.57	240	735760	40.00	ug/l	-0.04
44) Perylene-d12	37.79	264	492238	40.00	ug/l	-0.02
System Monitoring Compounds						
28) TCMX	23.23	209	174103	34.29	ug/L	0.00
Spiked Amount 40.000			Recovery = 85.72%			
Target Compounds						Qvalue
2) N-nitrosodimethylamine	0.00	74	0	N.D.		
3) bis(2-Chloroethyl)ether	0.00	93	0	N.D.		
4) 1,3-Dichlorobenzene	12.14	146	128	N.D.		
5) 1,4-Dichlorobenzene	12.14	146	128	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-Nitrosodi-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	22.56	149	444	N.D.		
26) 4-Chlorophenylphenylether	23.23	204	1025	N.D.		
27) Fluorene	0.00	166	0	N.D.		
29) Azobenzene	0.00	77	0	N.D.		
31) N-Nitrosodiphenylamine	23.23	168	302	N.D.		
32) 4-Bromophenylphenylether	0.00	248	0	N.D.		
33) Hexachlorobenzene	0.00	284	0	N.D.		
34) Phenanthrene	25.51	178	727	N.D.		

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

Quantitation Report (QT Reviewed)

Data File : C:\HPCHEM\1\DATA\MAY2402\12154.D Vial: 8
Acq On : 24 May 2002 10:26 pm Operator: Morgan
Sample : gc/ms scan 5/22 fb Inst : 209
Misc : Multiplr: 1.00
MS Integration Params: GOOFY.P
Quant Time: May 28 11:47 2002 Quant Results File: GCMS0520.RES

Quant Method : C:\HPCHEM\1\METHODS\GCMS0520.M (RTE Integrator)
Title : 209 625/TCLP calibration table
Last Update : Tue May 28 11:39:13 2002
Response via : Initial Calibration
DataAcq Meth : GCMS0520

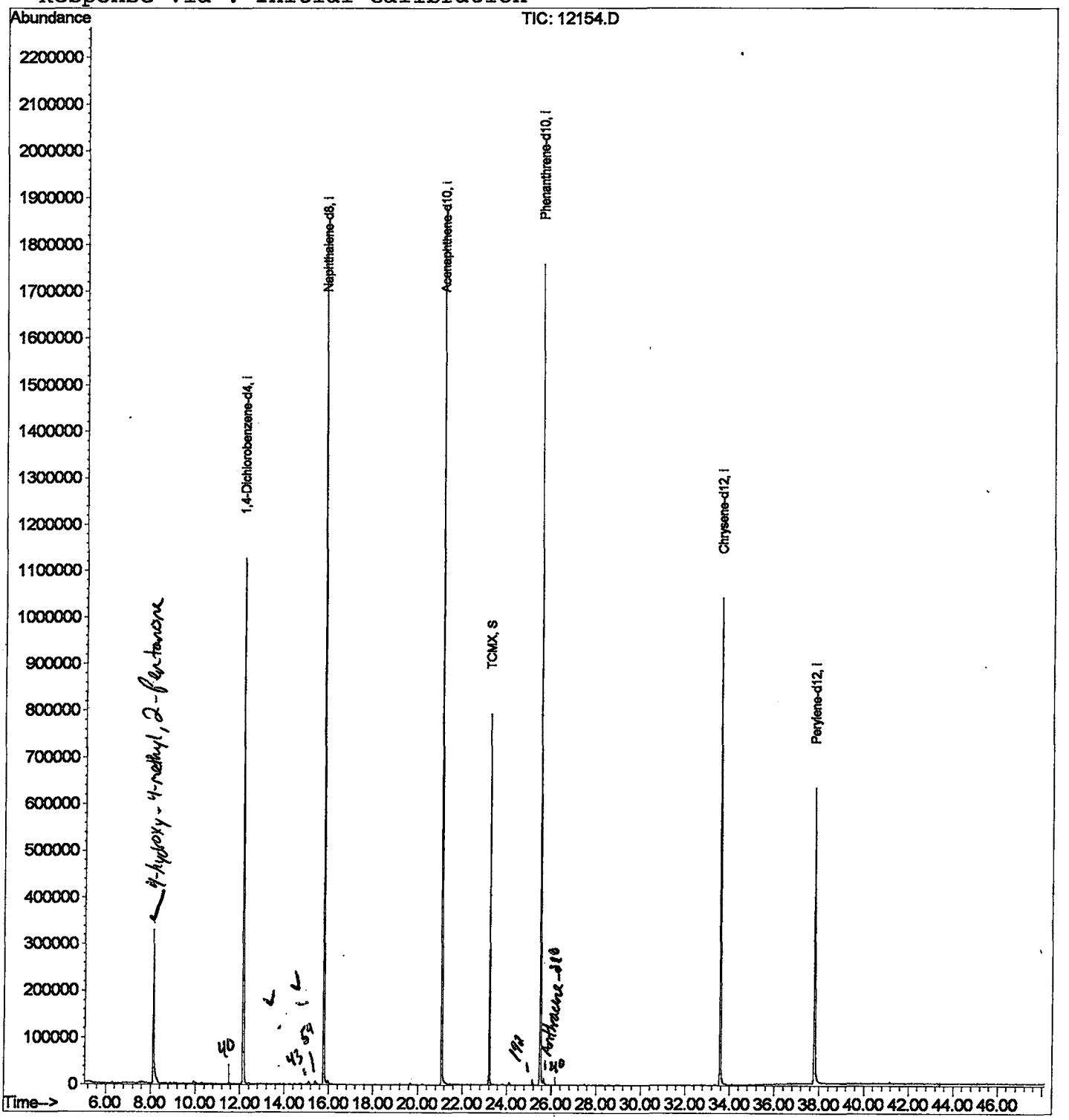
Compound	R.T.	QIon	Response	Conc Unit	Qvalue
35) Anthracene	25.51	178	727	N.D.	
36) Di-n-butylphthalate	27.49	149	3228	N.D.	
37) Fluoranthene	0.00	202	0	N.D.	
39) Pyrene	0.00	202	0	N.D.	
40) Butylbenzylphthalate	0.00	149	0	N.D.	
41) Benzo(a)anthracene	33.57	228	1976	N.D.	
42) Bis(2-ethylhexyl)phthalate	33.80	149	545	N.D.	
43) Chrysene	33.64	228	215	N.D.	
45) Di-n-Octylphthalate	0.00	149	0	N.D.	
46) Benzo(b)fluoranthene	0.00	252	0	N.D.	
47) Benzo(k)fluoranthene	0.00	252	0	N.D.	
48) Benzo(a)pyrene	37.78	252	2019	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\MAY2402\12154.D Vial: 8
Acq On : 24 May 2002 10:26 pm Operator: Morgan
Sample : gc/ms scan 5/22 fb Inst : 209
Misc : Multiplr: 1.00
MS Integration Params: GOOFY.P
Quant Time: May 28 11:47 2002 Quant Results File: GCMS0520.RES

Method : C:\HPCHEM\1\METHODS\GCMS0520.M (RTE Integrator)
Title : 209 625/TCLP calibration table
Last Update : Tue May 28 11:39:13 2002
Response via : Initial Calibration



LSC Area Percent Report

Data File : C:\HPCHEM\1\DATA\MAY2402\12154.D Vial: 8
Acq On : 24 May 2002 10:26 pm Operator: Morgan
Sample : gc/ms scan 5/22 fb Inst : 209
Misc : Multiplr: 1.00
MS Integration Params: LSCINT.P

Method : C:\HPCHEM\1\METHODS\GCMS0520.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

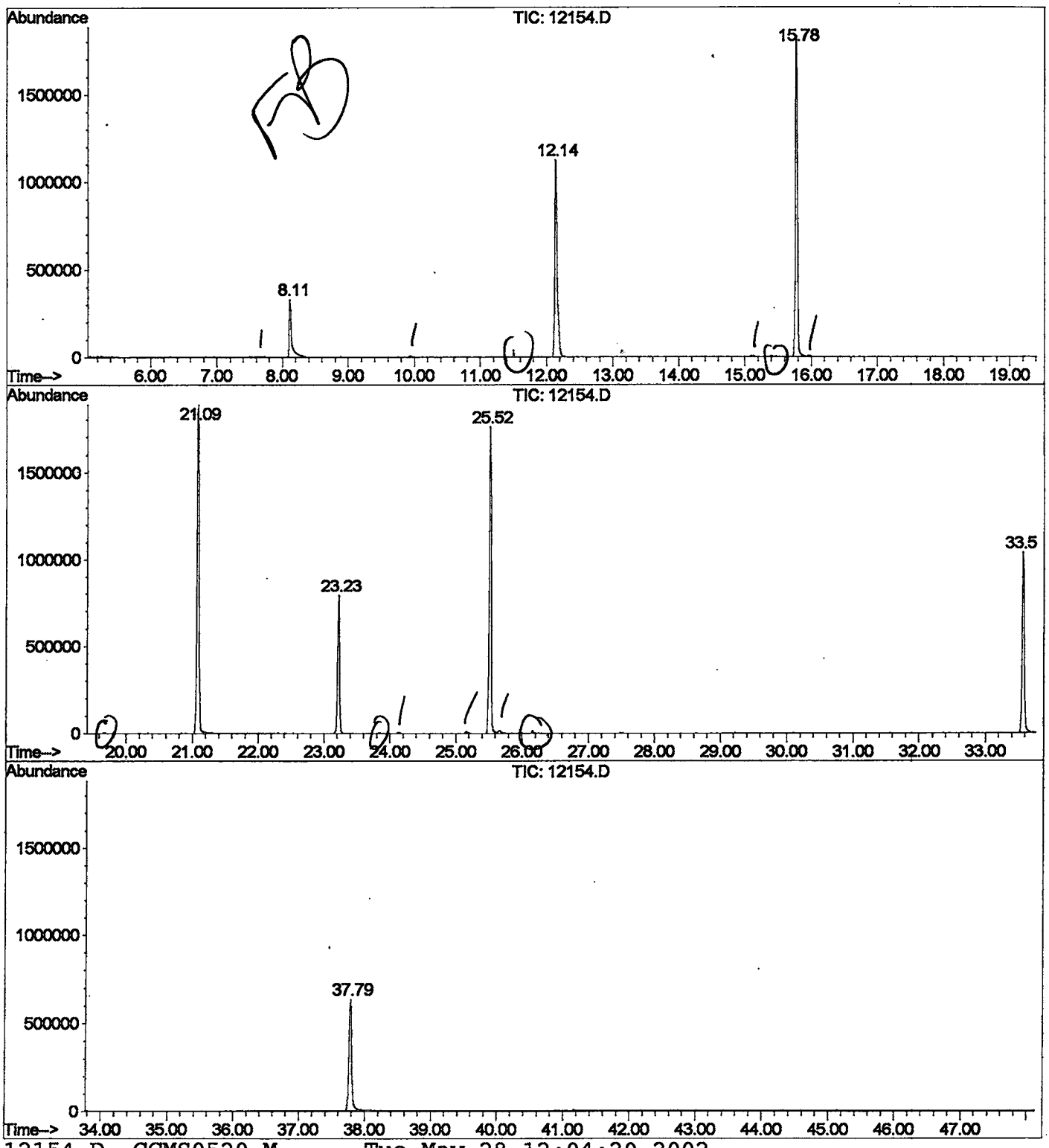
peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	peak area	peak % max.	% of total
1	8.114	330	334	361	rBV	328606	818887	20.52%	3.874%
2	12.135	767	772	786	rBV	1126127	2879233	72.15%	13.621%
3	15.780	1163	1169	1186	rBV	1831938	3902646	97.80%	18.463%
4	21.086	1740	1747	1762	rBV	1885763	3990508	100.00%	18.879%
5	23.225	1971	1980	1988	rBV	792449	1620961	40.62%	7.669%
6	25.520	2223	2230	2241	rBV2	1758153	3721052	93.25%	17.604%
7	33.571	3101	3107	3127	rBV2	1041168	2343032	58.72%	11.085%
8	37.785	3558	3566	3590	rBV	630856	1861241	46.64%	8.805%

Sum of corrected areas: 21137560

12154.D GCMS0520.M Tue May 28 12:04:19 2002

LSC Report - Integrated Chromatogram

File : C:\HPCHEM\1\DATA\MAY2402\12154.D
Operator : Morgan
Acquired : 24 May 2002 10:26 pm using AcqMethod GCMS0520
Instrument : 209
Sample Name: gc/ms .scan 5/22 fb
Misc Info :
Vial Number: 8
Quant File :GCMS0520.RES (RTE Integrator)



12154.D GCMS0520.M Tue May 28 12:04:20 2002

Library Search Compound Report

Data File : C:\HPCHEM\1\DATA\MAY2402\12154.D
Acq On : 24 May 2002 10:26 pm
Sample : gc/ms scan 5/22 fb
Misc :
MS Integration Params: LSCINT.P

Vial: 8
Operator: Morgan
Inst : 209
Multiplr: 1.00

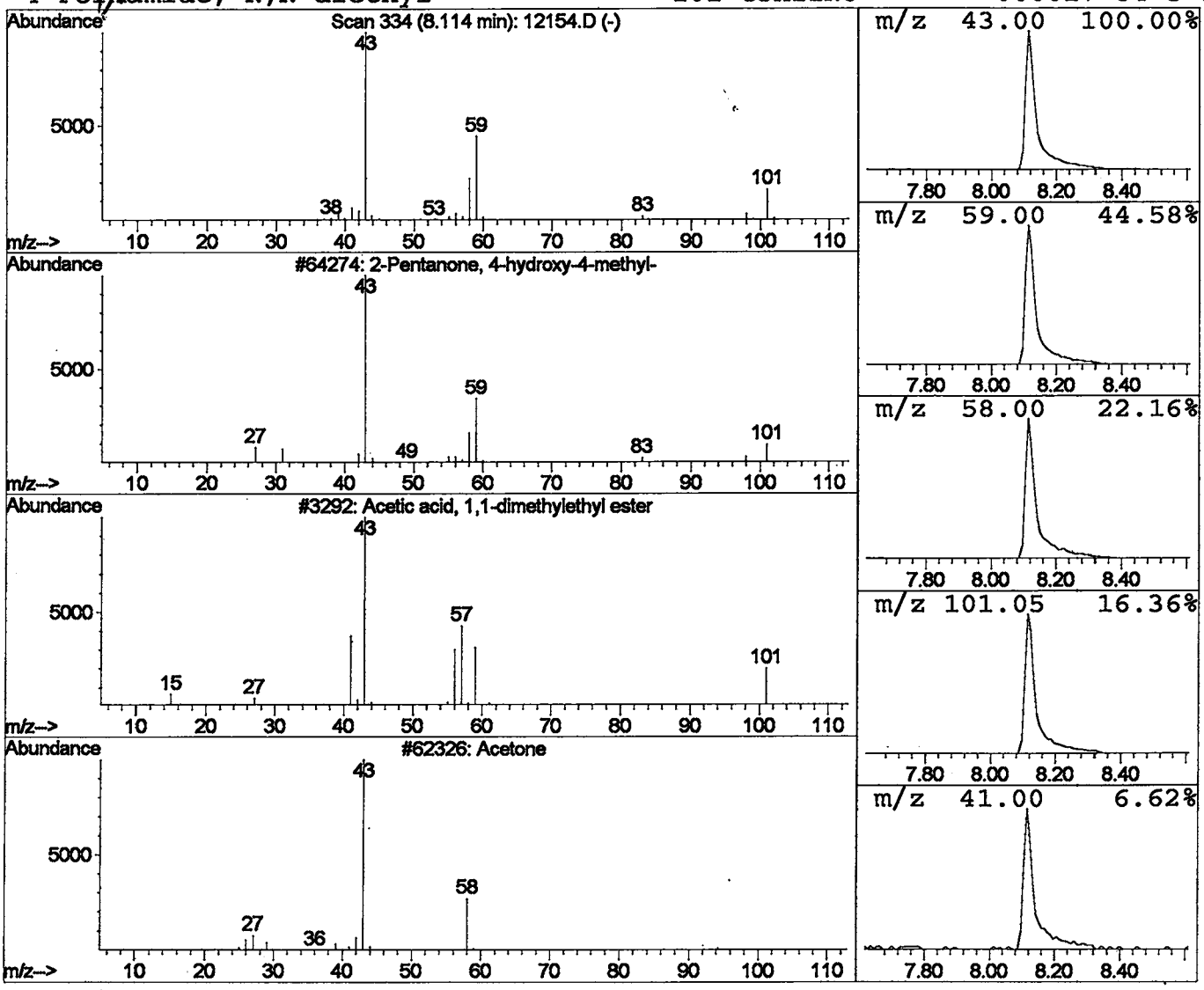
Quant Method : C:\HPCHEM\1\METHODS\GCMS0520.M (RTE Integrator)
Title : 209 625/TCLP calibration table
Library : C:\DATABASE\NBS75K.L

****

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.11	11.38 ug/l	818887	1,4-Dichlorobenzene-d4	12.14

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	83
2	Acetic acid, 1,1-dimethylethyl este	116	C6H12O2	000540-88-5	25
3	Acetone	58	C3H6O	000067-64-1	16
4	Formamide, N,N-diethyl-	101	C5H11NO	000617-84-5	12



Tentatively Identified Compound (LSC) summary

Operator ID: Morgan Date Acquired: 24 May 2002 10:26 pm
Data File: C:\HPCHEM\1\DATA\MAY2402\12154.D
Name: gc/ms scan 5/22 fb
Misc:
Method: C:\HPCHEM\1\METHODS\GCMS0520.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	8.11	11.4	ug/l	818887	ISTD01	12.14	2879230	40.0

12154.D GCMS0520.M Tue May 28 12:04:21 2002