User: Vinci, David L
Westchester Dept. of Labs & Research
Date: 06/07/2002 Time: 11:16:57

Sample: AE12679
Analysis: $GCMS GCMS Scan For Unknowns
Units: ug
Started: 6/7/02 11:16 Ended: 6/7/02 11:16 Analyst: DLV
Page: 1

	Component Name	Viol	Result	Qualifier	MDL
1	Other unknown compounds		.		
2	Surr_2,4-DDD		75		10.0

Comments for Sample AE12679 - Analysis $GCMS

Westchester Dept. of Labs & Research
User: Vinci, David L
Date: 06-07-2002 Time: 11:17:29

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\JUN0302\12679.D Vial: 11
Acq On : 3 Jun 2002 7:03 pm Operator:
Sample : gc/ms scan 5/31 fb Inst : 209
Misc : Multiplr: 1.00
MS Integration Params: GOOFY.P
Quant Time: Jun 4 9:00 2002 Quant Results File: GCMS0531.RES

Quant Method : C:\HPCHEM\1\METHODS\GCMS0531.M (RTE Integrator)
Title : 209 625/TCLP calibration table
Last Update : Fri May 31 09:43:55 2002
Response via : Initial Calibration
DataAcq Meth : GCMS0531

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	12.07	152	475388	40.00	ug/l	-0.01
10) Naphthalene-d8	15.69	136	1888434	40.00	ug/l	-0.02
17) Acenaphthene-d10	20.99	164	889830	40.00	ug/l	-0.01
30) Phenanthrene-d10	25.42	188	1370955	40.00	ug/l	-0.01
38) Chrysene-d12	33.48	240	781969	40.00	ug/l	-0.03
44) Perylene-d12	37.65	264	467860	40.00	ug/l	-0.03
System Monitoring Compounds						
28) TCMX	23.13	209	184333	29.94	ug/L	-0.02
Spiked Amount 40.000			Recovery = 74.85%			
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	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-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.	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	22.50	149	857	N.D.		
26) 4-Chlorophenylphenylether	23.14	204	1191	N.D.		
27) Fluorene	0.00	166	0	N.D.		
29) Azobenzene	0.00	77	0	N.D.		
31) N-Nitrosodiphenylamine	23.13	168	614	N.D.		
32) 4-Bromophenylphenylether	0.00	248	0	N.D.		
33) Hexachlorobenzene	0.00	284	0	N.D.		
34) Phenanthrene	25.43	178	532	N.D.		

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

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Misc : Multiplr: 1.00
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Last Update : Fri May 31 09:43:55 2002
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DataAcq Meth : GCMS0531

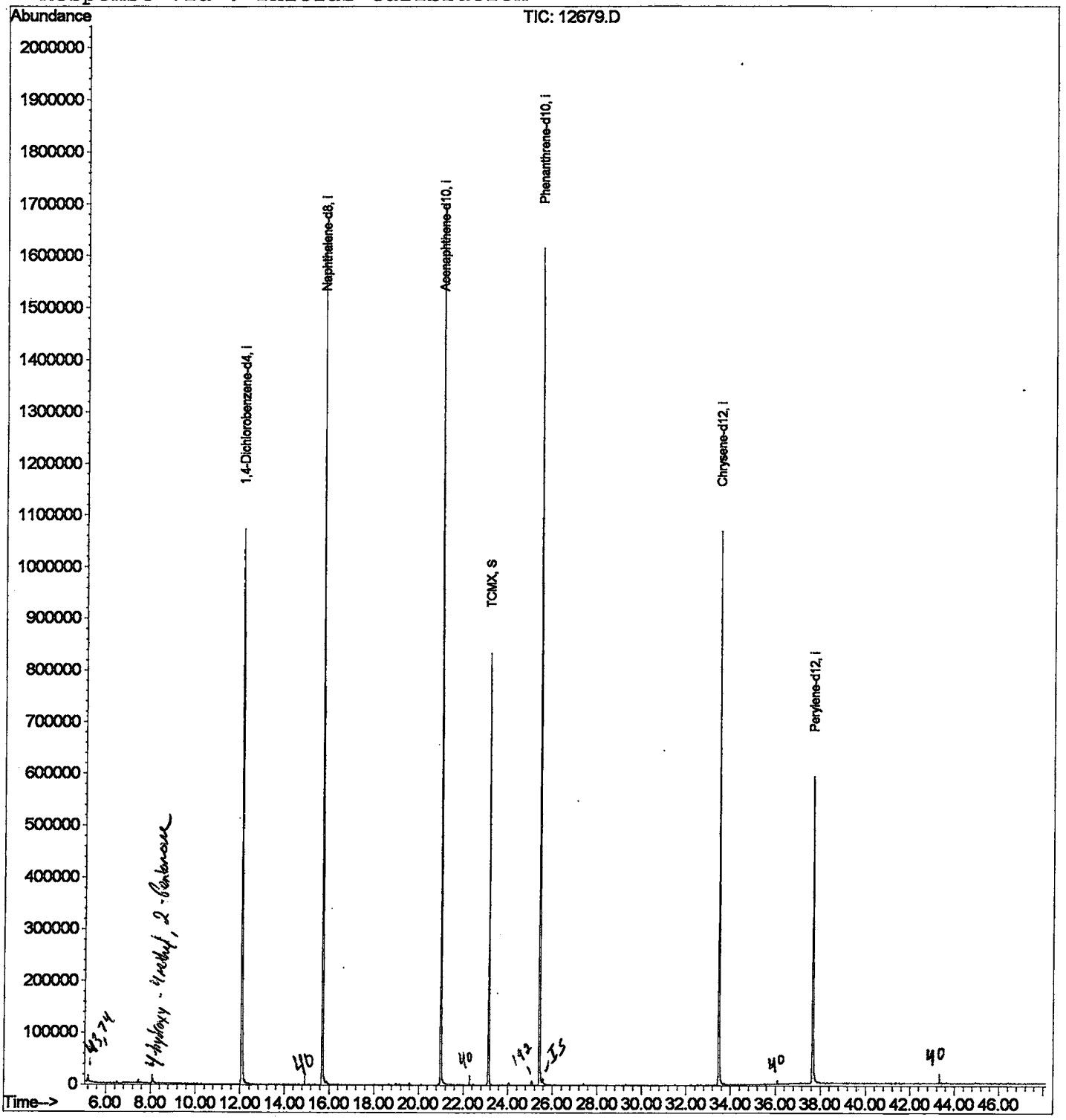
	Compound	R.T.	QIon	Response	Conc Unit	Qvalue
35)	Anthracene	25.43	178	532	N.D.	
36)	Di-n-butylphthalate	27.42	149	4928	N.D.	
37)	Fluoranthene	0.00	202	0	N.D.	
39)	Pyrene	0.00	202	0	N.D.	
40)	Butylbenzylphthalate	0.00	149	0	N.D.	
41)	Benzo(a)anthracene	33.48	228	1837	N.D.	
42)	Bis(2-ethylhexyl)phthalate	33.71	149	1157	N.D.	
43)	Chrysene	33.48	228	1837	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	37.65	252	1830	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\JUN0302\12679.D Vial: 11
Acq On : 3 Jun 2002 7:03 pm Operator:
Sample : gc/ms scan 5/31 fb Inst : 209
Misc : Multiplr: 1.00
MS Integration Params: GOOFY.P
Quant Time: Jun 4 9:00 2002 Quant Results File: GCMS0531.RES

Method : C:\HPCHEM\1\METHODS\GCMS0531.M (RTE Integrator)
Title : 209 625/TCLP calibration table
Last Update : Fri May 31 09:43:55 2002
Response via : Initial Calibration



LSC Area Percent Report

Data File : C:\HPCHEM\1\DATA\JUN0302\12679.D Vial: 11
Acq On : 3 Jun 2002 7:03 pm Operator:
Sample : gc/ms scan 5/31 fb Inst : 209
Misc : Multiplr: 1.00
MS Integration Params: LSCINT.P

Method : C:\HPCHEM\1\METHODS\GCMS0531.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

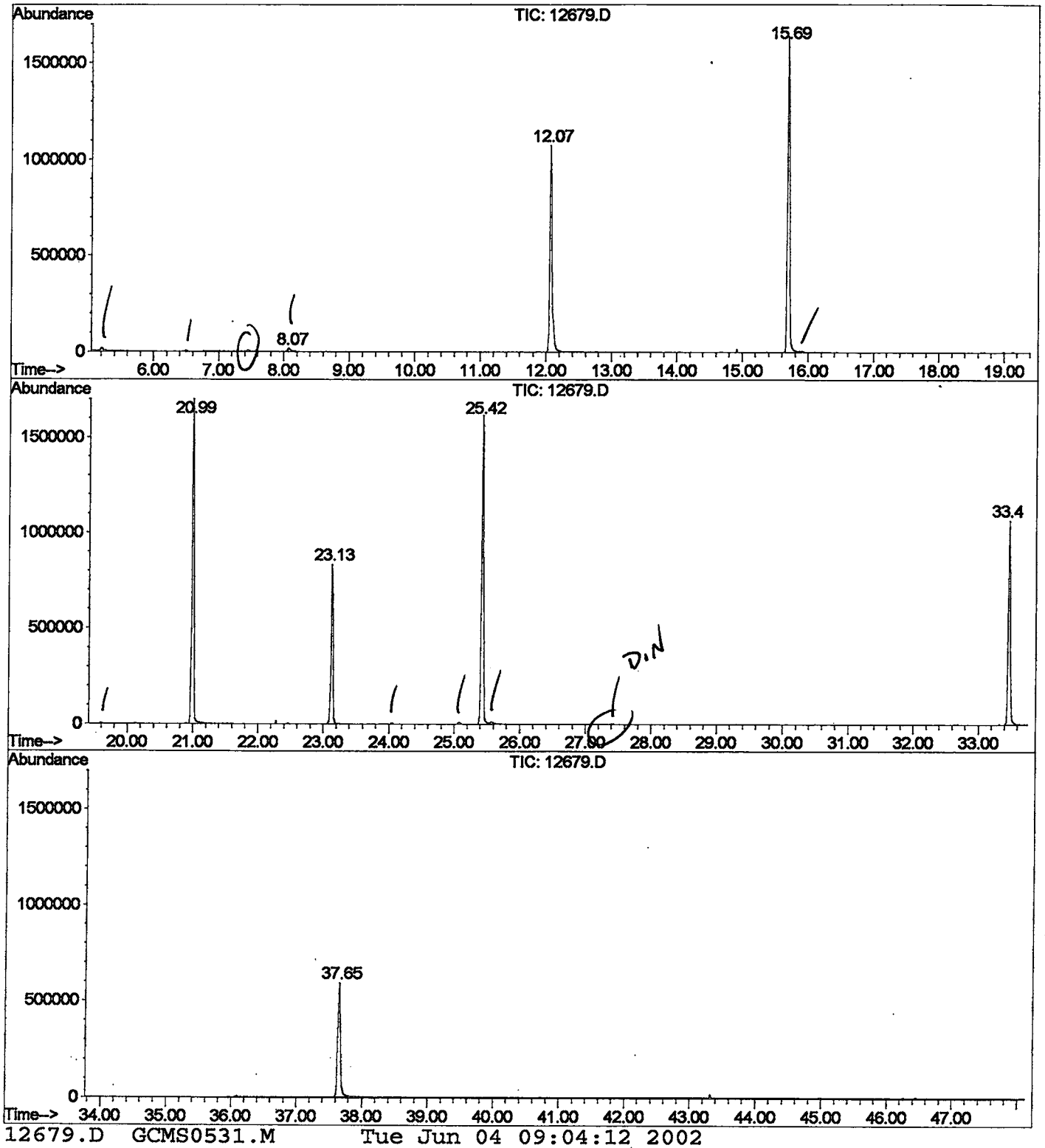
peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	peak area	peak % max.	% of total
1	8.073	326	330	336	rBV	16336	37694	1.06%	0.197%
2	12.067	760	765	785	rBV	1072921	2590502	72.97%	13.556%
3	15.693	1154	1160	1175	rBV	1634641	3489054	98.28%	18.258%
4	20.990	1731	1737	1755	rBV	1700715	3549993	100.00%	18.577%
5	23.129	1961	1970	1979	rBV	833918	1736202	48.91%	9.086%
6	25.424	2213	2220	2232	rBV2	1614836	3528616	99.40%	18.465%
7	33.475	3090	3097	3112	rBV2	1068844	2444402	68.86%	12.792%
8	37.652	3545	3552	3568	rBV2	590805	1732971	48.82%	9.069%

Sum of corrected areas: 19109434

12679.D GCMS0531.M Tue Jun 04 09:04:11 2002

LSC Report - Integrated Chromatogram

File : C:\HPCHEM\1\DATA\JUN0302\12679.D
Operator :
Acquired : 3 Jun 2002 7:03 pm using AcqMethod GCMS0531
Instrument : 209
Sample Name: gc/ms scan 5/31 fb
Misc Info :
Vial Number: 11
Quant File :GCMS0531.RES (RTE Integrator)



12679.D GCMS0531.M Tue Jun 04 09:04:12 2002

Library Search Compound Report

Data File : C:\HPCHEM\1\DATA\JUN0302\12679.D Vial: 11
Acq On : 3 Jun 2002 7:03 pm Operator:
Sample : gc/ms scan 5/31 fb Inst : 209
Misc : Multiplr: 1.00
MS Integration Params: LSCINT.P

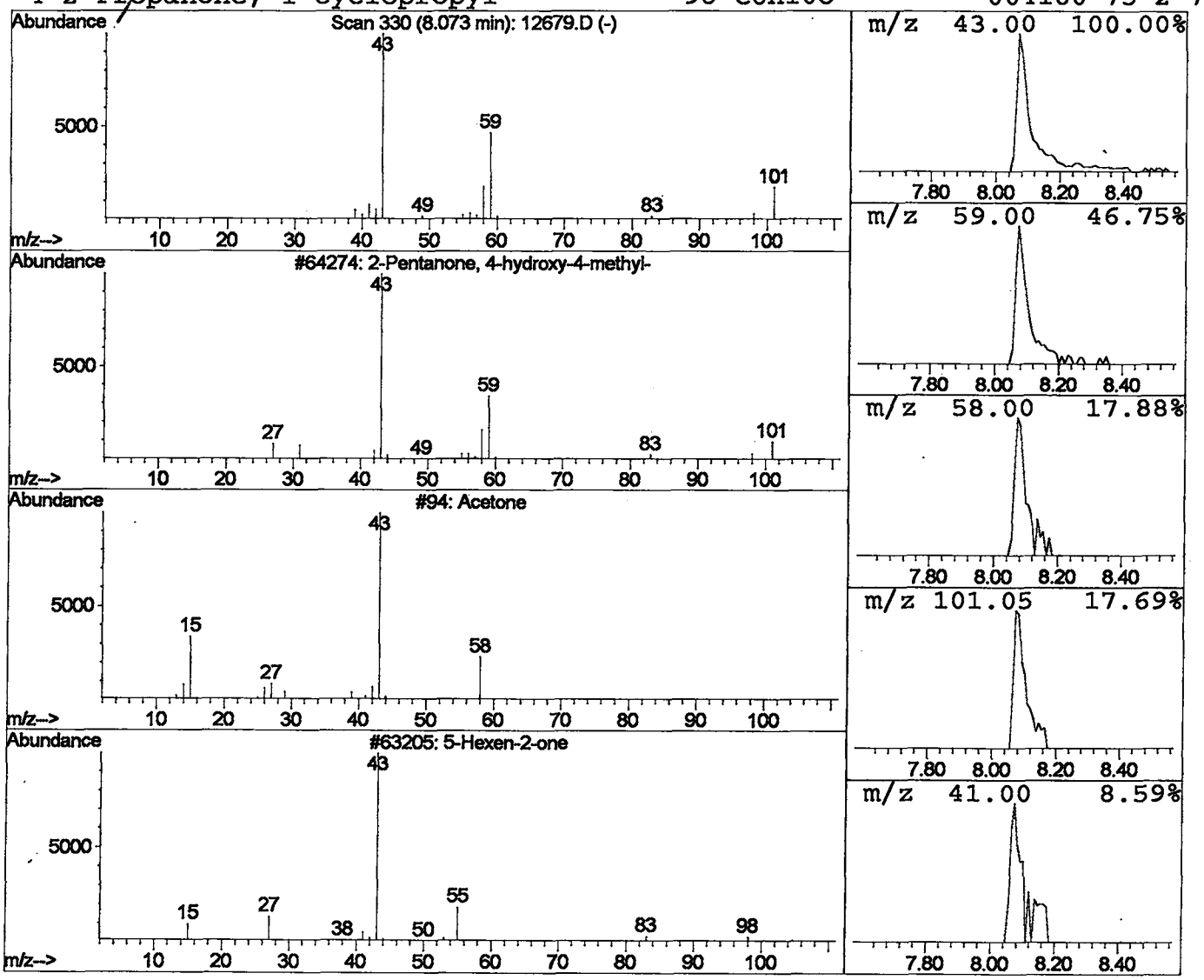
Quant Method : C:\HPCHEM\1\METHODS\GCMS0531.M (RTE Integrator)
Title : 209 625/TCLP calibration table
Library : C:\DATABASE\NBS75K.L

****

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.07	0.58 ug/l	37694	1,4-Dichlorobenzene-d4	12.07

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	83
2	Acetone	58	C3H6O	000067-64-1	7
3	5-Hexen-2-one	98	C6H10O	000109-49-9	7
4	2-Propanone, 1-cyclopropyl-	98	C6H10O	004160-75-2	7



Tentatively Identified Compound (LSC) summary

Operator ID: Date Acquired: 3 Jun 2002 7:03 pm
Data File: C:\HPCHEM\1\DATA\JUN0302\12679.D
Name: gc/ms scan 5/31 fb
Misc:
Method: C:\HPCHEM\1\METHODS\GCMS0531.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	8.07	0.6	ug/l	37694	ISTD01	12.07	2590500	40.0

12679.D GCMS0531.M Tue Jun 04 09:04:13 2002