User: Galos, Marie
Westchester Dept. of Labs & Research
Date: 06/03/2002 Time: 10:37:20

Sample: AE12098
Analysis: $B1GCMS Blank for GCMS Scan
Units: ug
Started: 5/31/20 00:02 Ended: 5/31/20 00:02 Analyst: MG
Result source: manual entry
Page: 1

	Component Name	Viol	Result	MDL
1	n-Nitrosodimethylamine		< MRL	2.0
2	bis(2-Chloroethyl)ether		< MRL	2.0
3	1,3-Dichlorobenzene		< MRL	2.0
4	1,4-Dichlorobenzene		< MRL	2.0
5	1,2-Dichlorobenzene		< MRL	2.0
6	2,2'-oxybis(1-Chloropropane)		< MRL	2.0
7	n-Nitrosodi-n-propylamine		< MRL	2.0
8	Hexachloroethane		< MRL	2.0
9	Nitrobenzene		< MRL	2.0
10	Isophorone		< MRL	2.0
11	bis(2-Chloroethoxy)methane		< MRL	2.0
12	1,2,4-Trichlorobenzene		< MRL	2.0
13	Naphthalene		< MRL	2.0
14	Hexachlorobutadiene		< MRL	2.0
15	2-Chloronaphthalene		< MRL	2.0
16	Dimethylphthalate		< MRL	2.0
17	2,6-Dinitrotoluene		< MRL	2.0
18	Acenaphthylene		< MRL	2.0
19	Acenaphthene		< MRL	2.0
20	2,4-Dinitrotoluene		< MRL	2.0
21	Diethylphthalate		< MRL	2.0
22	4-Chlorophenylphenylether		< MRL	2.0
23	Fluorene		< MRL	2.0
24	Azobenzene		< MRL	2.0
25	n-Nitrosodiphenylamine		< MRL	2.0
26	4-Bromophenylphenylether		< MRL	2.0
27	Hexachlorobenzene		< MRL	2.0
28	Phenanthrene		< MRL	2.0
29	Anthracene		< MRL	2.0
30	Di-n-butylphthalate		< MRL	2.0
31	Fluoranthene		< MRL	2.0
32	Pyrene		< MRL	2.0
33	Butylbenzylphthalate		< MRL	2.0
34	Benzo(a)anthracene		< MRL	2.0
35	bis(2-Ethylhexyl)phthalate		< MRL	2.0
36	Chrysene		< MRL	2.0
37	Di-n-octylphthalate		< MRL	2.0
38	Benzo(b)fluoranthene		< MRL	2.0
39	Benzo(k)fluoranthene		< MRL	2.0
40	Benzo(a)pyrene		< MRL	2.0
41	Indeno(1,2,3-cd)pyrene		< MRL	2.0
42	Dibenz(a,h)anthracene		< MRL	2.0
43	Benzo(g,h,i)perylene		< MRL	2.0

Quantitation Report (QT Reviewed)

Data File : C:\HPCHEM\1\DATA\MAY3102\LB.D Vial: 3
Acq On : 31 May 2002 1:54 pm Operator: Morgan
Sample : gc/ms scan 5/29 Inst : 209
Misc : Multiplr: 1.00
MS Integration Params: GOOFY.P
Quant Time: Jun 3 9:22 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	464616	40.00	ug/l	-0.01
10) Naphthalene-d8	15.69	136	1880290	40.00	ug/l	-0.02
17) Acenaphthene-d10	20.99	164	887383	40.00	ug/l	-0.01
30) Phenanthrene-d10	25.42	188	1408330	40.00	ug/l	-0.01
38) Chrysene-d12	33.47	240	876536	40.00	ug/l	-0.03
44) Perylene-d12	37.66	264	563796	40.00	ug/l	-0.02
System Monitoring Compounds						
28) TCMX	23.14	209	213021	34.70	ug/L	-0.01
Spiked Amount 40.000			Recovery = 86.75%			
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	11.97	146	573	N.D.		
5) 1,4-Dichlorobenzene	12.11	146	462	N.D.		
6) 1,2-Dichlorobenzene	12.63	146	185	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	13.48	117	290	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	15.58	180	208	N.D.		
15) Naphthalene	15.75	128	728	N.D.		
16) Hexachlorobutadiene	16.32	225	724	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	662	N.D.		
26) 4-Chlorophenylphenylether	23.13	204	1462	N.D.		
27) Fluorene	0.00	166	0	N.D.		
29) Azobenzene	0.00	77	0	N.D.		
31) N-Nitrosodiphenylamine	23.14	168	694	N.D.		
32) 4-Bromophenylphenylether	0.00	248	0	N.D.		
33) Hexachlorobenzene	0.00	284	0	N.D.		
34) Phenanthrene	25.43	178	495	N.D.		

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

Quantitation Report (QT Reviewed)

Data File : C:\HPCHEM\1\DATA\MAY3102\LB.D Vial: 3
Acq On : 31 May 2002 1:54 pm Operator: Morgan
Sample : gc/ms scan 5/29 Inst : 209
Misc : Multiplr: 1.00
MS Integration Params: GOOFY.P
Quant Time: Jun 3 9:22 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

	Compound	R.T.	QIon	Response	Conc Unit	Qvalue
35)	Anthracene	25.43	178	495	N.D.	
36)	Di-n-butylphthalate	27.41	149	73044	1.44 ug/l	99
37)	Fluoranthene	0.00	202	0	N.D.	
39)	Pyrene	0.00	202	0	N.D.	
40)	Butylbenzylphthalate	31.92	149	427	N.D.	
41)	Benzo(a)anthracene	33.48	228	3134	N.D.	
42)	Bis(2-ethylhexyl)phthalate	33.70	149	1832	N.D.	
43)	Chrysene	33.55	228	1169	N.D.	
45)	Di-n-Octylphthalate	0.00	149	0	N.D.	
46)	Benzo(b)fluoranthene	36.59	252	2046	N.D.	
47)	Benzo(k)fluoranthene	36.59	252	1270	N.D.	
48)	Benzo(a)pyrene	37.47	252	576	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

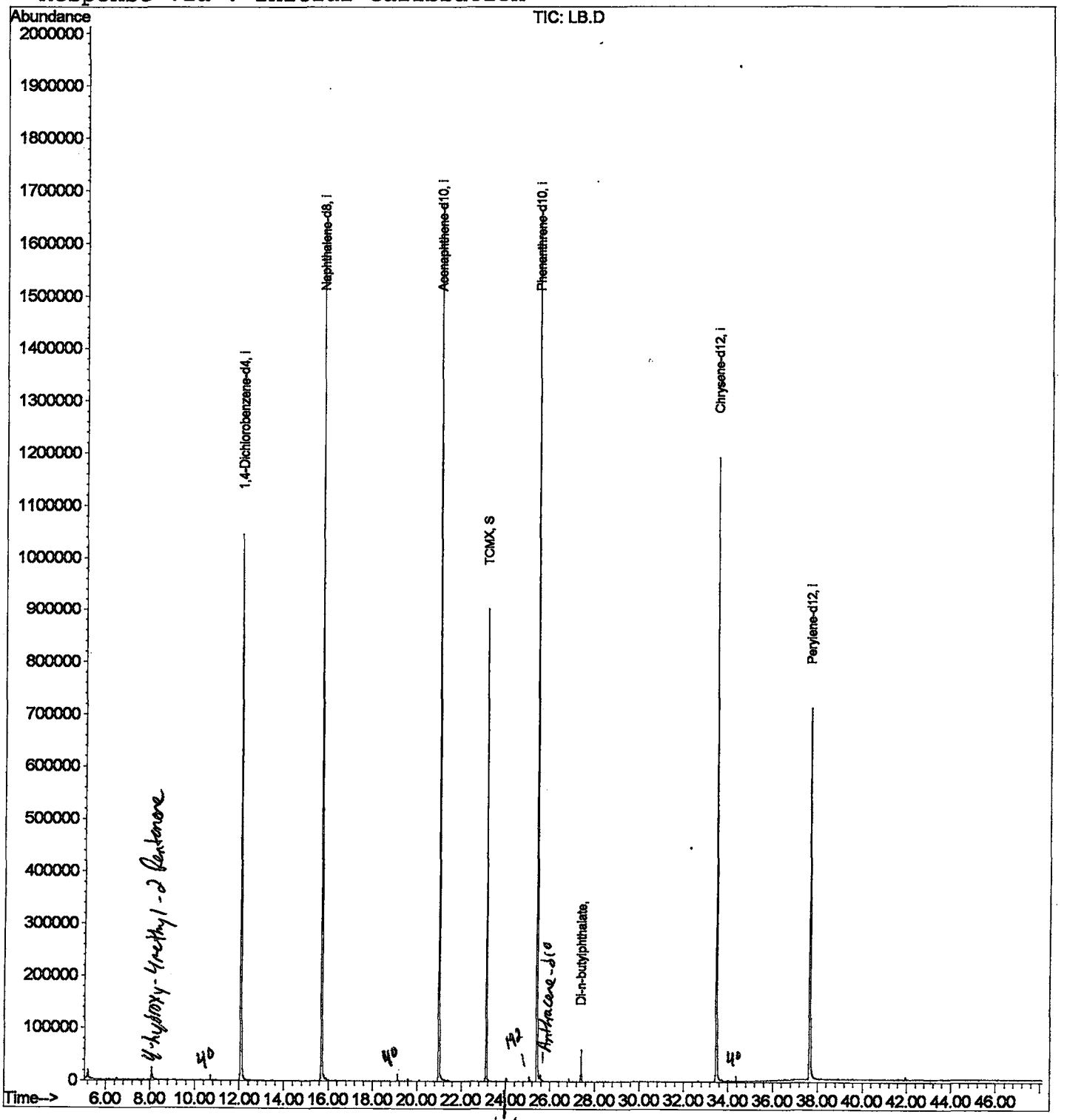
Quantitation Report

Data File : C:\HPCHEM\1\DATA\MAY3102\LB.D
Acq On : 31 May 2002 1:54 pm
Sample : gc/ms scan 5/29
Misc :
MS Integration Params: GOOFY.P
Quant Time: Jun 3 9:22 2002

Vial: 3
Operator: Morgan
Inst : 209
Multiplr: 1.00

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\MAY3102\LB.D Vial: 3
Acq On : 31 May 2002 1:54 pm Operator: Morgan
Sample : gc/ms scan 5/29 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

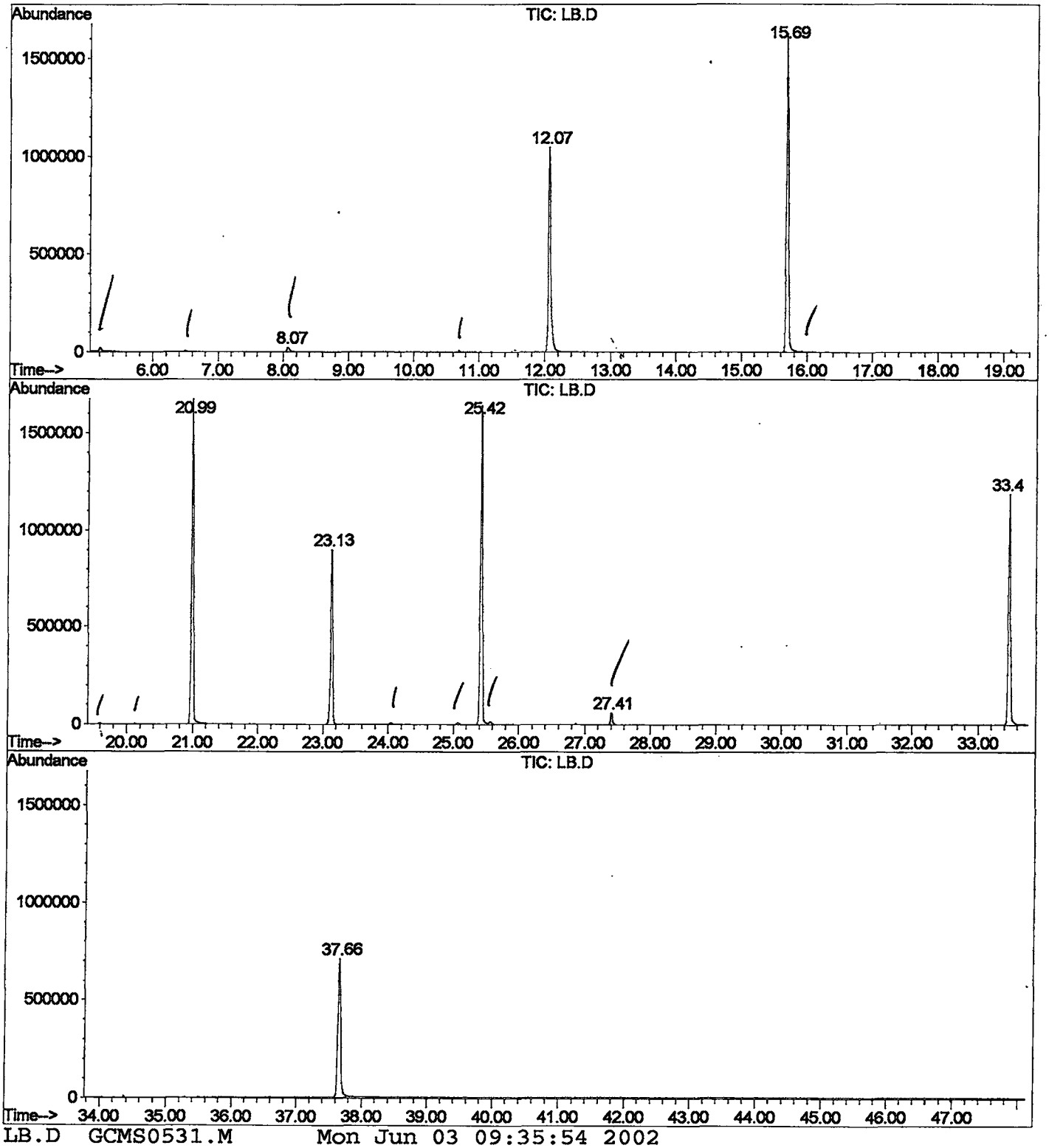
peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	peak area	peak % max.	% of total
1	8.072	326	330	340	rBV	23354	62666	1.74%	0.312%
2	12.065	760	765	784	rBV	1045038	2536168	70.59%	12.638%
3	15.692	1154	1160	1178	rBV	1627043	3448125	95.98%	17.183%
4	20.989	1730	1737	1751	rBV	1677059	3524351	98.10%	17.563%
5	23.128	1962	1970	1979	rBV	901924	1979718	55.10%	9.865%
6	25.423	2213	2220	2232	rBV2	1643652	3592639	100.00%	17.903%
7	27.406	2431	2436	2447	rVB	60470	117871	3.28%	0.587%
8	33.474	3089	3097	3113	rBV2	1192042	2739694	76.26%	13.653%
9	37.660	3544	3553	3565	rBV2	707269	2065969	57.51%	10.295%

Sum of corrected areas: 20067201

LB.D GCMS0531.M Mon Jun 03 09:35:53 2002

LSC Report - Integrated Chromatogram

File : C:\HPCHEM\1\DATA\MAY3102\LB.D
Operator : Morgan
Acquired : 31 May 2002 1:54 pm using AcqMethod GCMS0531
Instrument : 209
Sample Name: gc/ms scan 5/29
Misc Info :
Vial Number: 3
Quant File :GCMS0531.RES (RTE Integrator)



LB.D GCMS0531.M Mon Jun 03 09:35:54 2002

Library Search Compound Report

Data File : C:\HPCHEM\1\DATA\MAY3102\LB.D
Acq On : 31 May 2002 1:54 pm
Sample : gc/ms scan 5/29
Misc :
MS Integration Params: LSCINT.P

Vial: 3
Operator: Morgan
Inst : 209
Multiplr: 1.00

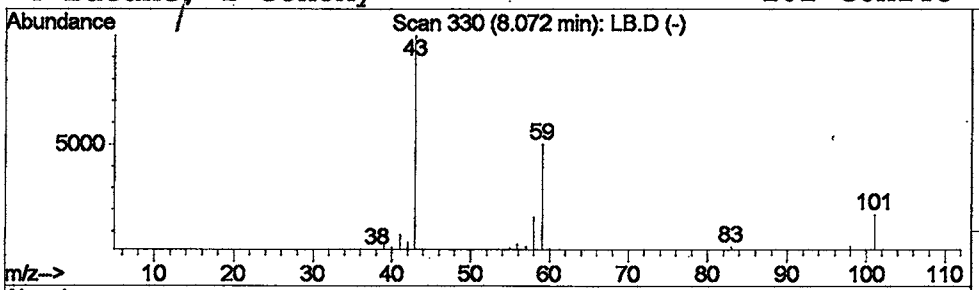
Quant Method : C:\HPCHEM\1\METHODS\GCMS0531.M (RTE Integrator)
Title : 209 625/TCLP calibration table
Library : C:\DATABASE\NBS75K.L

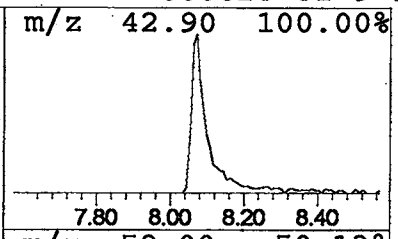
****

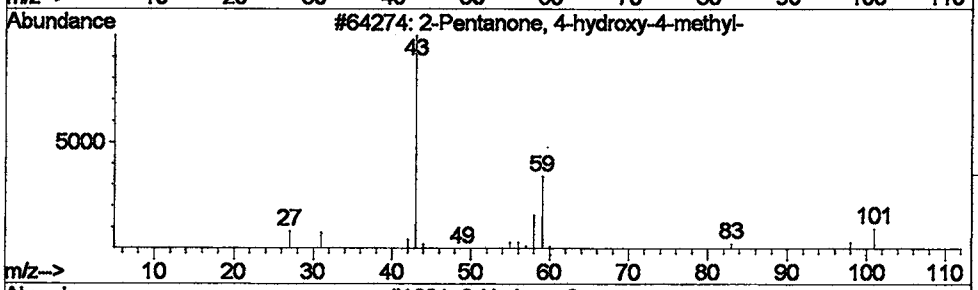
Peak Number 1 2-Pentanone, 4-hydroxy-4-methy Concentration Rank 1

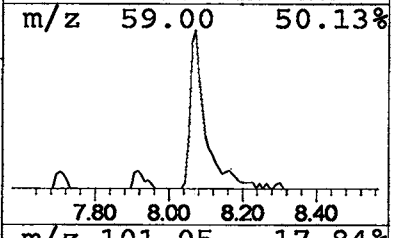
R.T.	EstConc	Area	Relative to ISTD	R.T.
8.07	0.99 ug/l	62666	1,4-Dichlorobenzene-d4	12.07

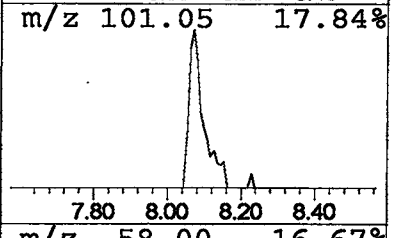
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	83
2	3-Hydroxy-2-pentanone	102	C5H10O2	003142-66-3	9
3	Morpholine, 4-methyl-	101	C5H11NO	000109-02-4	9
4	Butane, 1-ethoxy-	102	C6H14O	000628-81-9	9

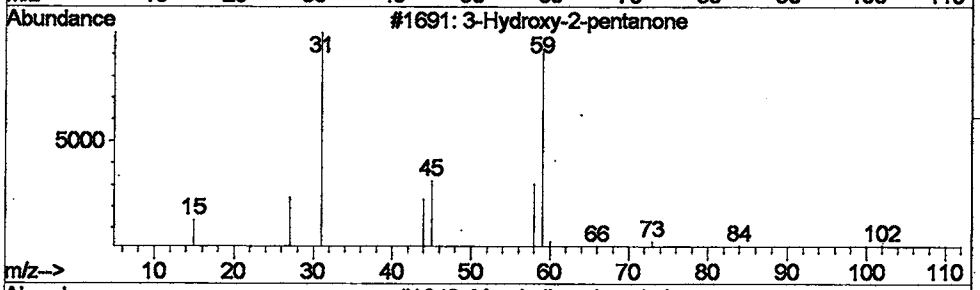


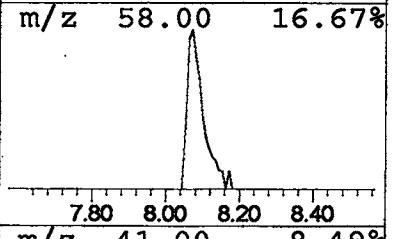


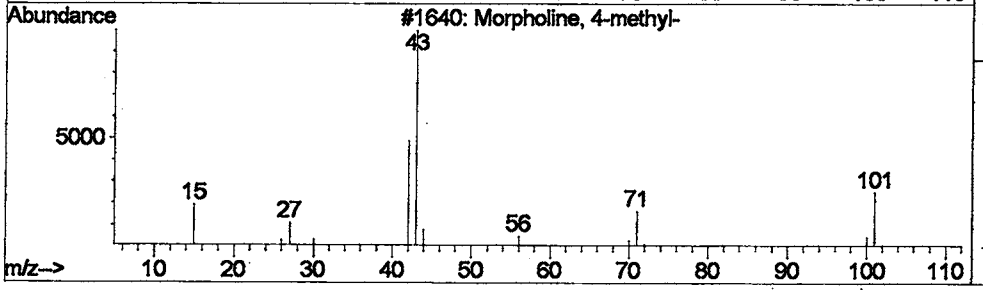


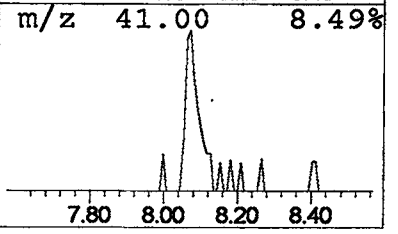












Tentatively Identified Compound (LSC) summary

Operator ID: Morgan Date Acquired: 31 May 2002 1:54 pm
Data File: C:\HPCHEM\1\DATA\MAY3102\LB.D
Name: gc/ms scan 5/29
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	1.0	ug/l	62666	ISTD01	12.07	2536170	40.0

LB.D GCMS0531.M Mon Jun 03 09:35:55 2002